LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143525.D
Acq On : 21 Aug 2025 12:46
Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 11M-N

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143525.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.275	6	14	26	rVB	805590	1313709	56.81%	4.778%
2	5.157	498	504	517	rBV	143607	222905	9.64%	0.811%
3	5.563	567	573	589	rBV	1589295	2097916	90.72%	7.629%
4	6.257	687	691	704	rBV2	10626	24786	1.07%	0.090%
5	6.557	736	742	758	rBV	1560605	2142617	92.65%	7.792%
6	6.922	799	804	812	rVB	541646	700686	30.30%	2.548%
7	7.481	893	899	911	rBV	1072438	1431019	61.88%	5.204%
8	8.204	1017	1022	1024	rBV	715965	955896	41.34%	3.476%
9	8.228	1024	1026	1032	rVV	655791	742357	32.10%	2.700%
10	8.281	1032	1035	1042	rVB	26905	37708	1.63%	0.137%
11	8.798	1119	1123	1132	rVB	39223	53823	2.33%	0.196%
12	8.922	1139	1144	1148	rBV	232753	298042	12.89%	1.084%
13	8.969	1148	1152	1156	rVV	149909	197886	8.56%	0.720%
14	9.016	1156	1160	1166	rVB	161138	210069	9.08%	0.764%
15	9.281	1199	1205	1210	rBV	1866336	2312517	100.00%	8.410%
16	9.381	1218	1222	1227	rVB	66221	86511	3.74%	0.315%
17	9.481	1231	1239	1245	rVV	19692	33255	1.44%	0.121%
18	9.545	1245	1250	1255	rVV	60744	97456	4.21%	0.354%
19	9.616	1255	1262	1265	rVV3	94593	188908	8.17%	0.687%
20	9.645	1265	1267	1275	rVB	46697	57532	2.49%	0.209%
21	9.733	1275	1282	1292	rBV3	34100	89410	3.87%	0.325%
22	9.822	1292	1297	1301	rVB2	23625	31674	1.37%	0.115%
23	9.963	1313	1321	1324	rBV	796547	1058770	45.78%	3.850%
24	9.992	1324	1326	1330	rVV	449727	511934	22.14%	1.862%
25	10.033	1330	1333	1336	rVB	24005	28783	1.24%	0.105%
26	10.116	1342	1347	1351	rVB3	17017	25197	1.09%	0.092%
27	10.169	1351	1356	1360	rBV	331411	429869	18.59%	1.563%
28	10.369	1386	1390	1397	rVB3	13617	24620	1.06%	0.090%
29	10.510	1409	1414	1419	rBV	510419	635315	27.47%	2.310%
30	10.592	1419	1428	1432	rVV2	45243	113894	4.93%	0.414%
31	10.669	1434	1441	1447	rVV	265465	362902	15.69%	1.320%
32	10.751	1450	1455	1470	rVB	1034352	1379832	59.67%	5.018%
33	10.875	1470	1476	1483	rBV	84034	121610	5.26%	0.442%
34	10.980	1491	1494	1500	rVB2	38858	49834	2.15%	0.181%
35	11.057	1500	1507	1510	rBV4	56155	101668	4.40%	0.370%
36	11.139	1517	1521	1524	rVB2	18790	24339	1.05%	0.089%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
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Acq On : 21 Aug 2025 12:46
Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.186	1524	1529	1531	rBV3	16306	27135	1.17%	0.099%
38	11.257	1537	1541	1545	rVB2	33836	42785	1.85%	0.156%
39	11.304	1545	1549	1552	rVV3	30380	46305	2.00%	0.168%
40	11.345	1552	1556	1562	rVB	92583	122567	5.30%	0.446%
41	11.451	1568	1574	1575	rBV	688391	793426	34.31%	2.885%
42	11.475	1575	1578	1583	rVV	1309686	1683242	72.79%	6.121%
43	11.527	1583	1587	1590	rVV	136446	189126	8.18%	0.688%
44	11.563	1590	1593	1598	rVB	48270	66446	2.87%	0.242%
45	11.680	1607	1613	1620	rBV	303877	408530	17.67%	1.486%
46	11.745	1620	1624	1628	rVV4	20566	36534	1.58%	0.133%
47	11.786	1628	1631	1636	rVV4	15018	27375	1.18%	0.100%
48	11.857	1637	1643	1650	rVB4	11952	27508	1.19%	0.100%
49	11.951	1654	1659	1661	rBV	75763	101599	4.39%	0.369%
50	11.980	1661	1664	1669	rVV	124768	164367	7.11%	0.598%
51	12.027	1669	1672	1674	rVV	26598	34112	1.48%	0.124%
52	12.063	1674	1678	1686	rVB2	147582	250377	10.83%	0.911%
53	12.245	1706	1709	1713	rBV	38086	42811	1.85%	0.156%
54	12.663	1776	1780	1784	rBV	466210	608747	26.32%	2.214%
55	12.698	1784	1786	1790	rVB3	56672	64746	2.80%	0.235%
56	12.851	1808	1812	1814	rBV	130512	159567	6.90%	0.580%
57	12.892	1814	1819	1833	rVB2	291984	519776	22.48%	1.890%
58	13.033	1836	1843	1851	rBV	1259684	1618667	70.00%	5.887%
59	13.116	1851	1857	1861	rBV3	20829	38462	1.66%	0.140%
60	13.221	1866	1875	1879	rBV3	56993	108678	4.70%	0.395%
61	13.292	1883	1887	1890	rVV	44750	55445	2.40%	0.202%
62	13.327	1890	1893	1901	rVB3	22181	40913	1.77%	0.149%
63	13.469	1912	1917	1922	rBV3	26952	43425	1.88%	0.158%
64	13.557	1927	1932	1936	rBV	76801	94828	4.10%	0.345%
65	14.092	2015	2023	2035	rVB2	488053	806377	34.87%	2.933%
66	14.863	2150	2154	2160	rVB	19410	28193	1.22%	0.103%
67	15.151	2198	2203	2211	rBV	58564	139636	6.04%	0.508%
68	15.210	2211	2213	2219	rVB2	24450	33380	1.44%	0.121%
69	15.463	2252	2256	2262	rVB	28276	43660	1.89%	0.159%
70	15.527	2262	2267	2272	rVV	32045	49321	2.13%	0.179%
71	15.592	2272	2278	2294	rVB	432576	721614	31.20%	2.624%
72	16.057	2352	2357	2363	rBV	17081	29957	1.30%	0.109%
73	17.127	2533	2539	2546	rBV2	13814	32587	1.41%	0.119%

Sum of corrected areas: 27497473

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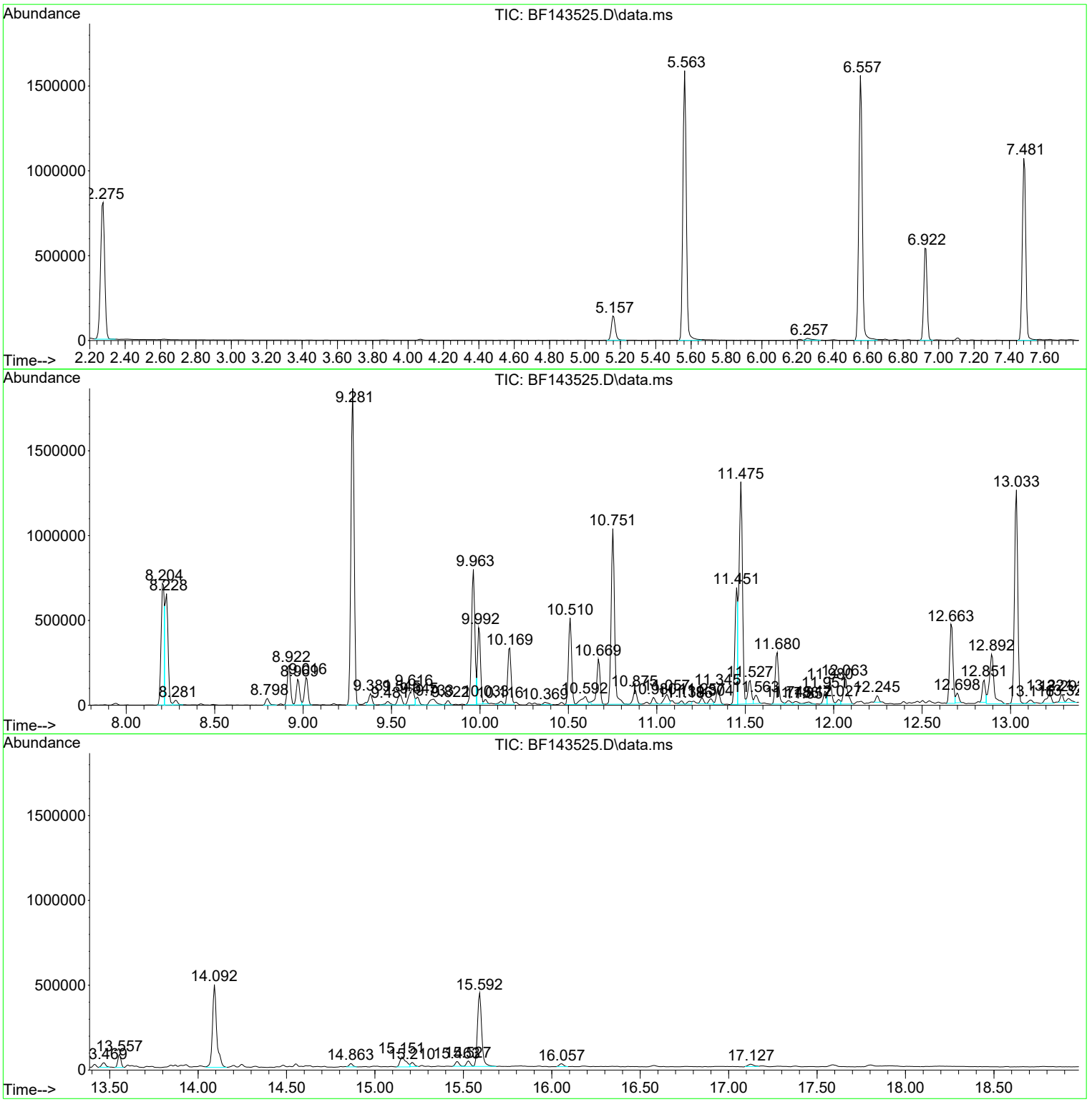
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

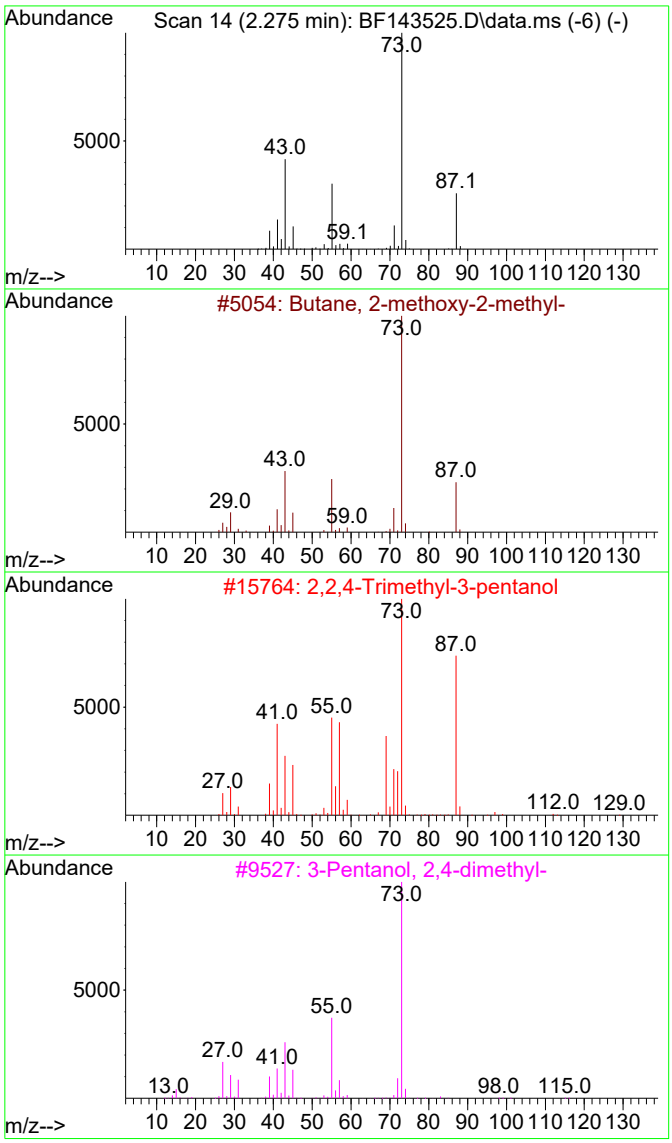
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

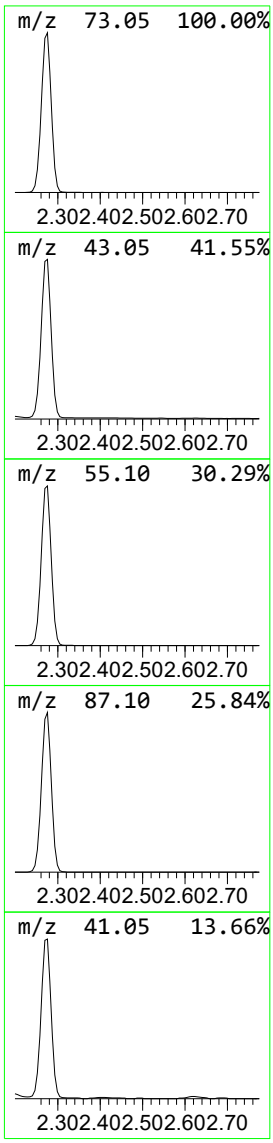
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.275	37.50 ng	1313710	1,4-Dichlorobenzene-d4	6.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143525.D
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Operator : RC/JU
Sample : Q2819-07
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

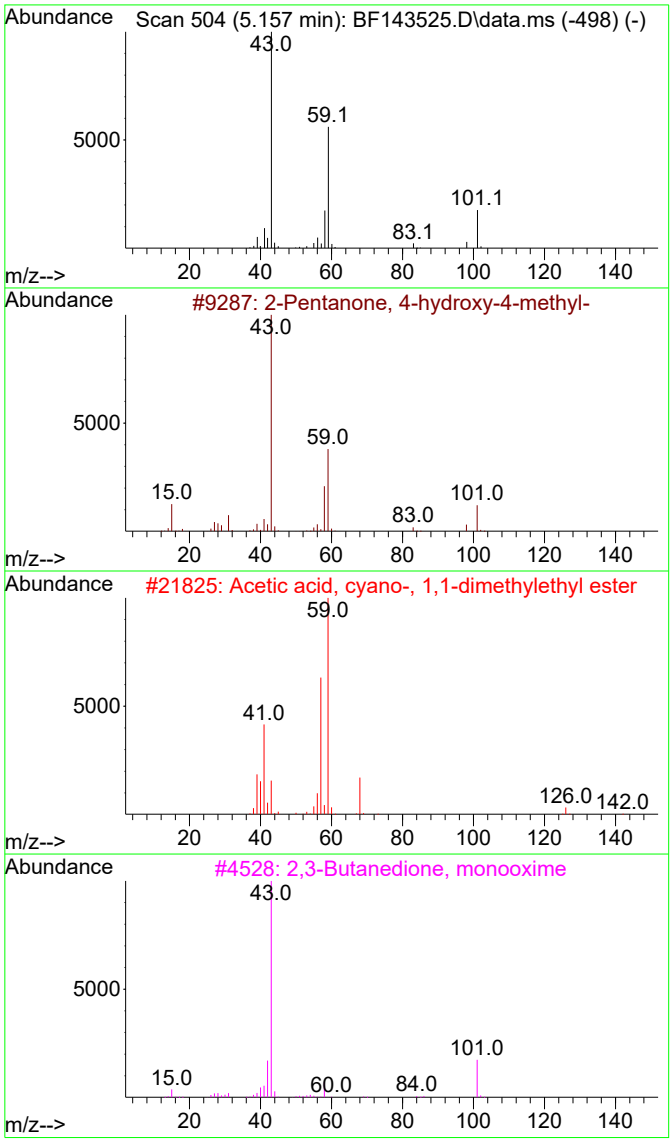
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

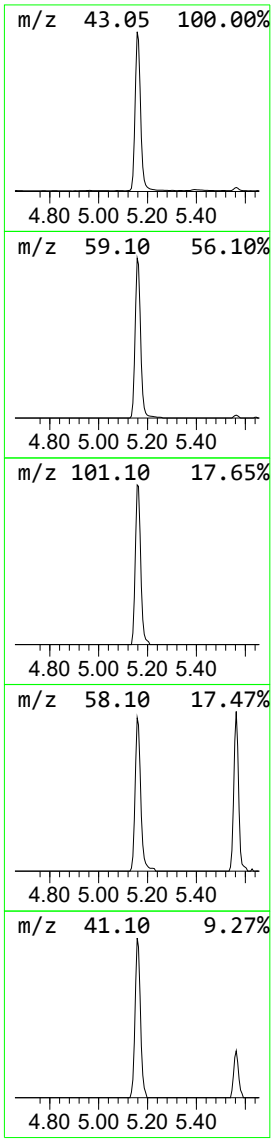
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	6.36 ng	222905	1,4-Dichlorobenzene-d4	6.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
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Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

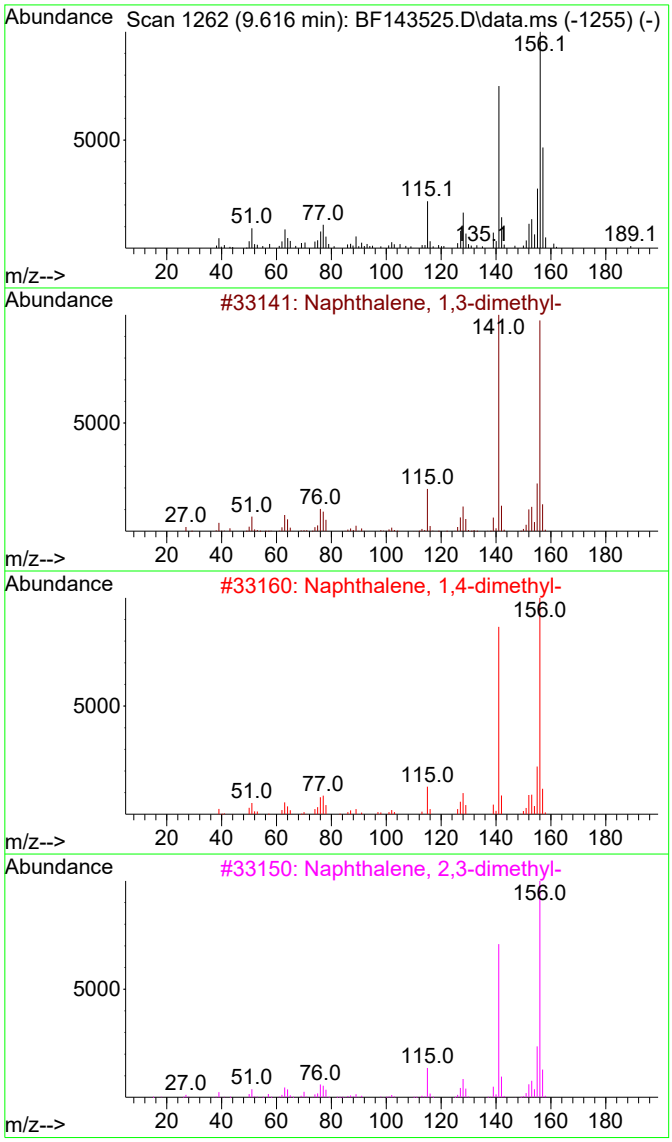
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

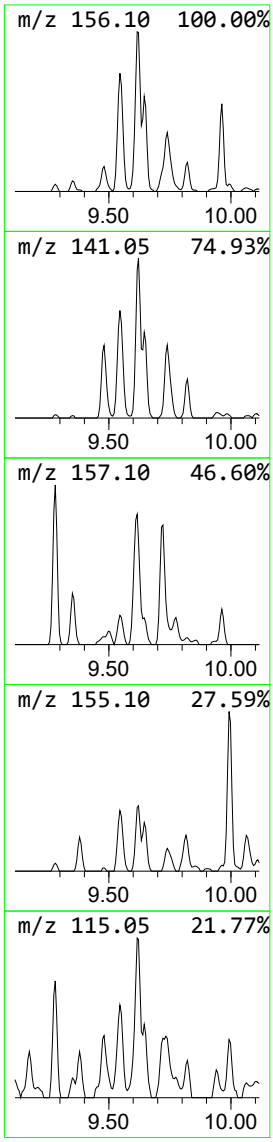
****

Peak Number 3 Naphthalene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.616	3.57 ng	188908	Acenaphthene-d10	9.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
2	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	94





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Sample : Q2819-07
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Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

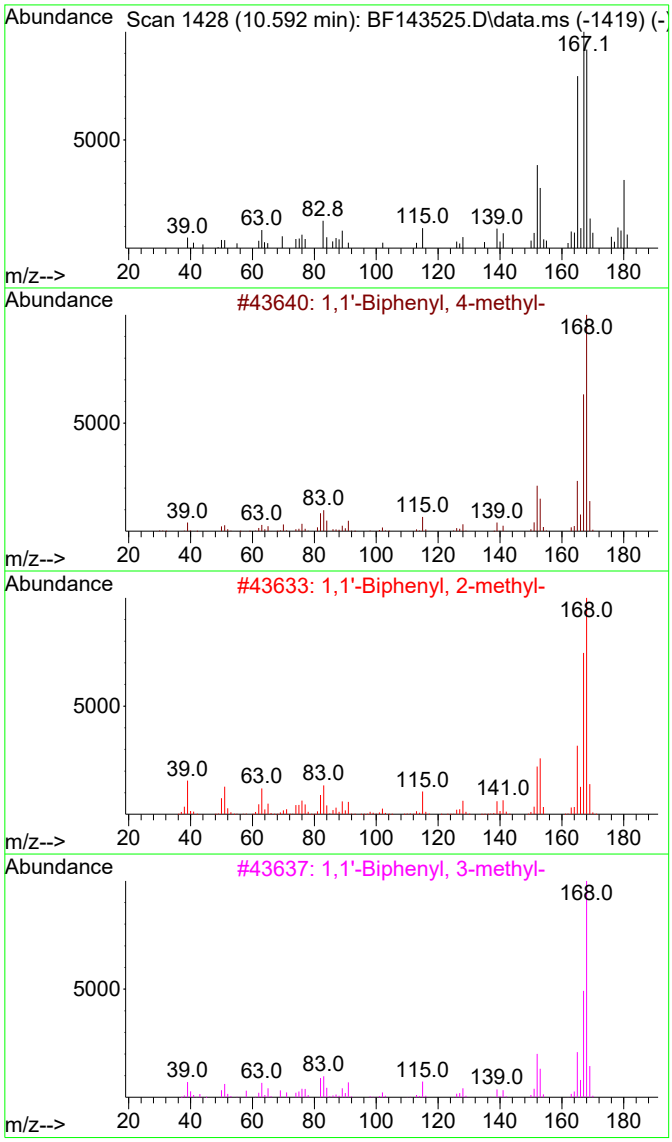
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

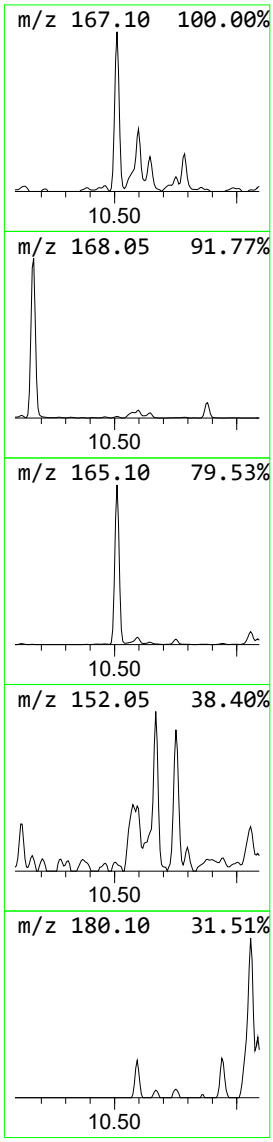
****

Peak Number 4 1,1'-Biphenyl, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.592	2.15 ng	113894	Acenaphthene-d10	9.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	87
2	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	83
3	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	70
4	Diphenylmethane	168	C13H12	000101-81-5	64
5	Fluorene, 1,4-dihydro-	168	C13H12	041593-21-9	60





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Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

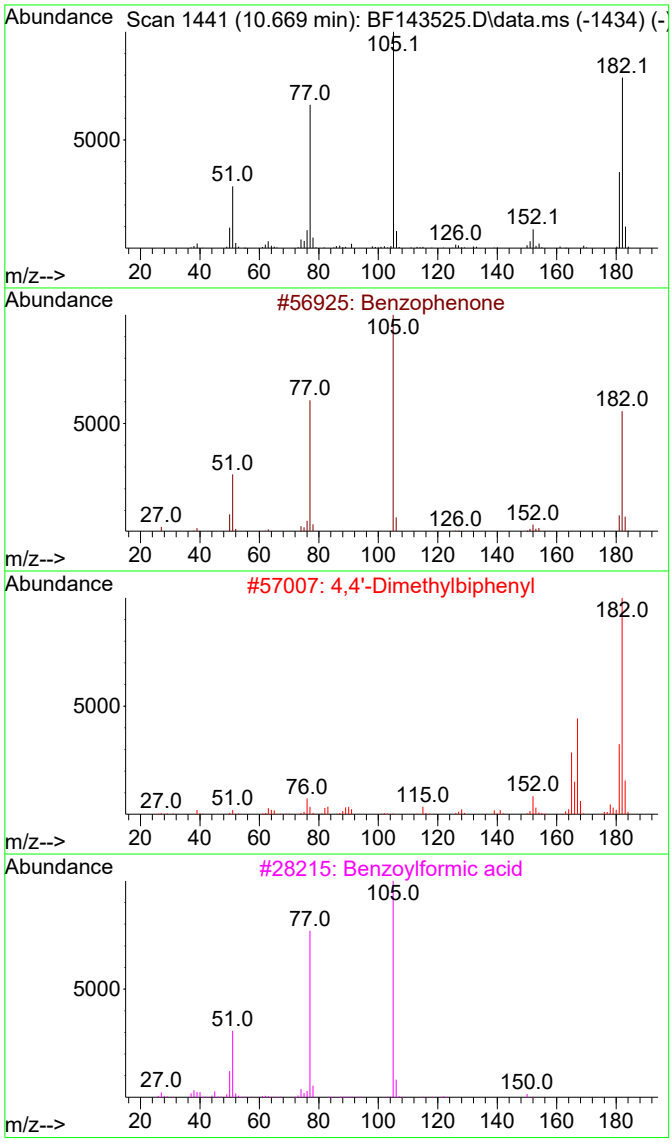
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

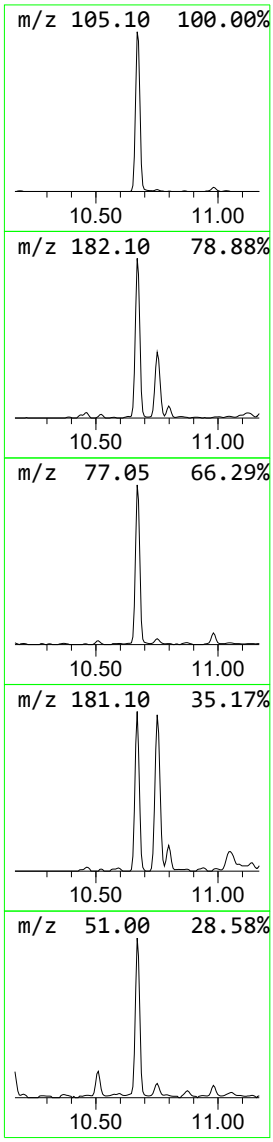
****

Peak Number 5 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.669	6.86 ng	362902	Acenaphthene-d10	9.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	94
2	4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	43
3	Benzoylformic acid	150	C8H6O3	000611-73-4	41
4	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	41
5	Ethanedione, diphenyl-	210	C14H10O2	000134-81-6	35





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

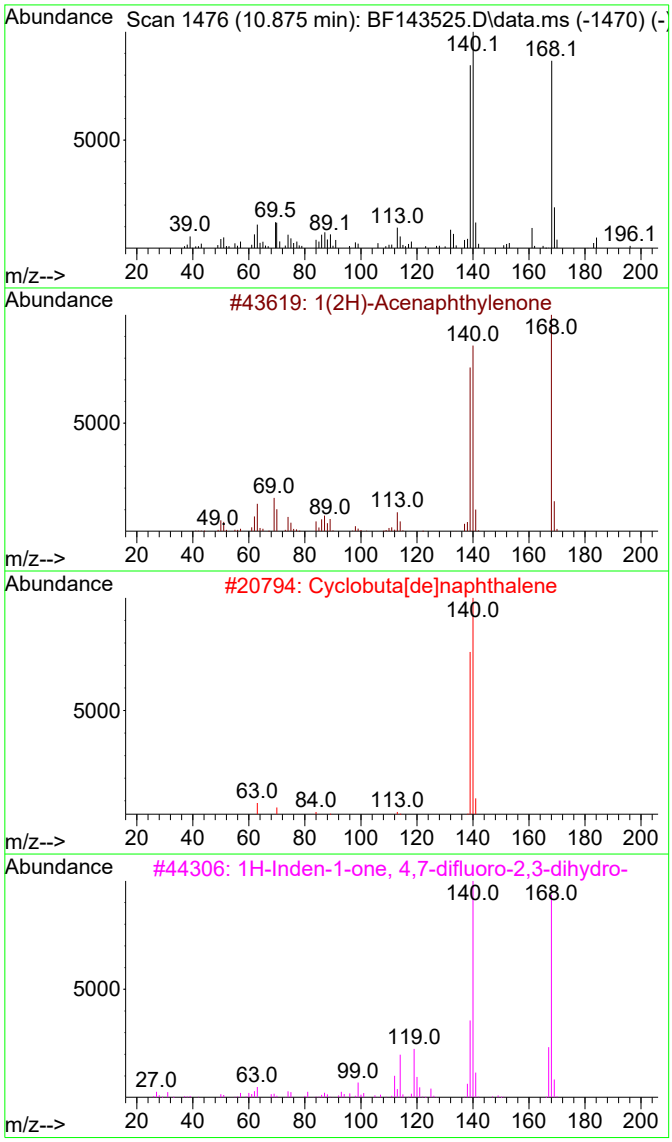
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

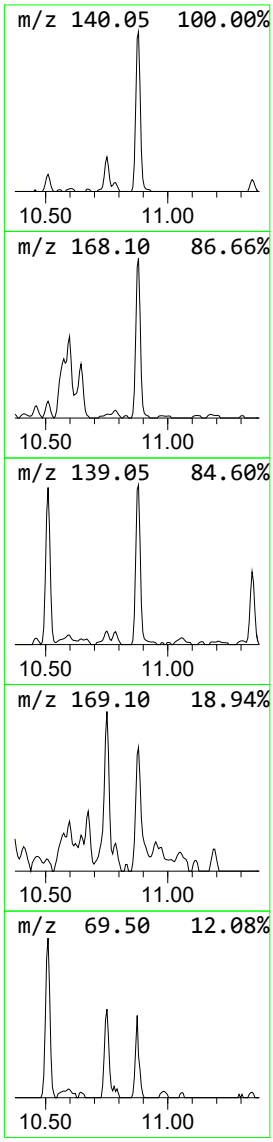
****

Peak Number 6 1(2H)-Acenaphthylenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.875	3.07 ng	121610	Phenanthrene-d10	11.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	95
2	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	60
3	1H-Inden-1-one, 4,7-difluoro-2,3...	168	C9H6F2O	130408-16-1	50
4	Thiophene, 2,5-dipropyl-	168	C10H16S	054411-07-3	49
5	Dibenzofuran	168	C12H8O	000132-64-9	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143525.D
Acq On : 21 Aug 2025 12:46
Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

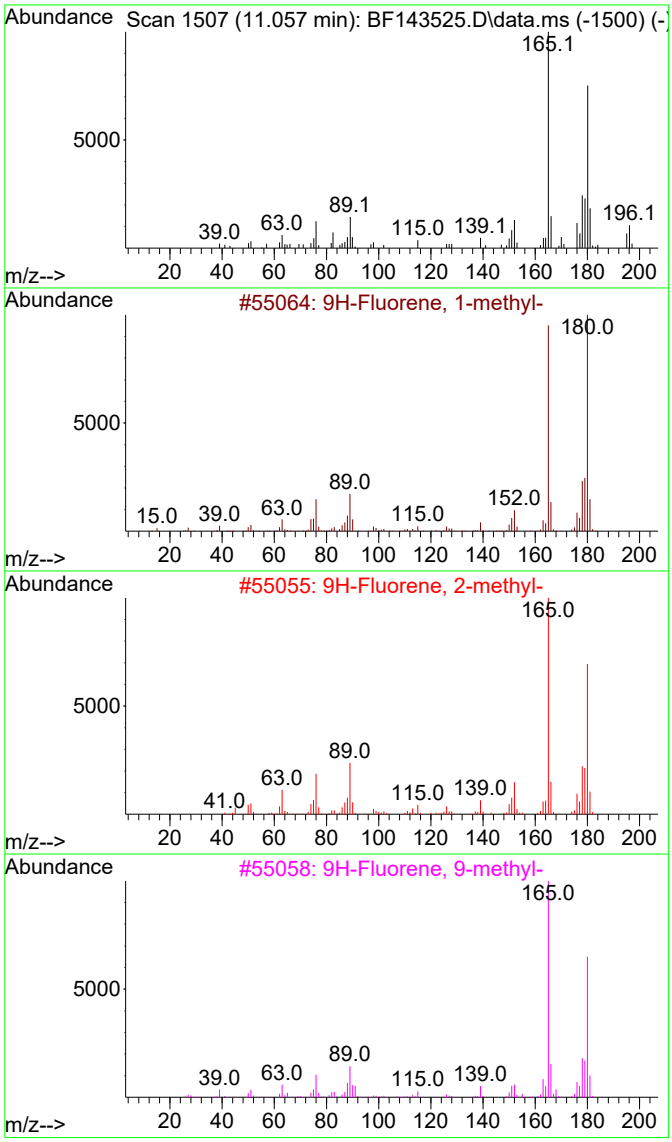
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

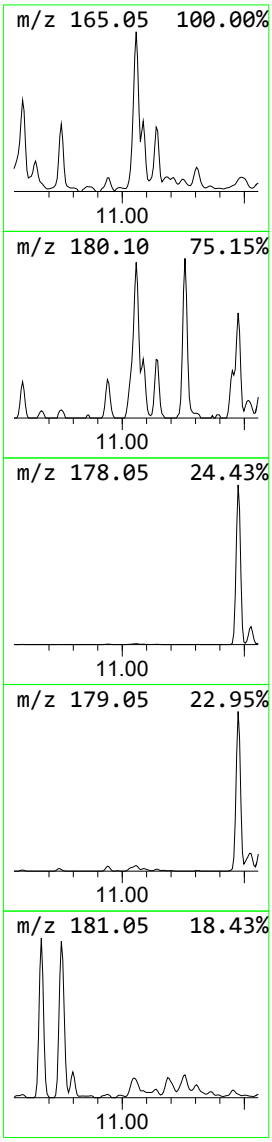
****

Peak Number 7 9H-Fluorene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.057	2.56 ng	101668	Phenanthrene-d10	11.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	96
2	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	96
3	9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	90
4	9H-Fluorene, 4-methyl-	180	C14H12	001556-99-6	90
5	9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	89





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Acq On : 21 Aug 2025 12:46
Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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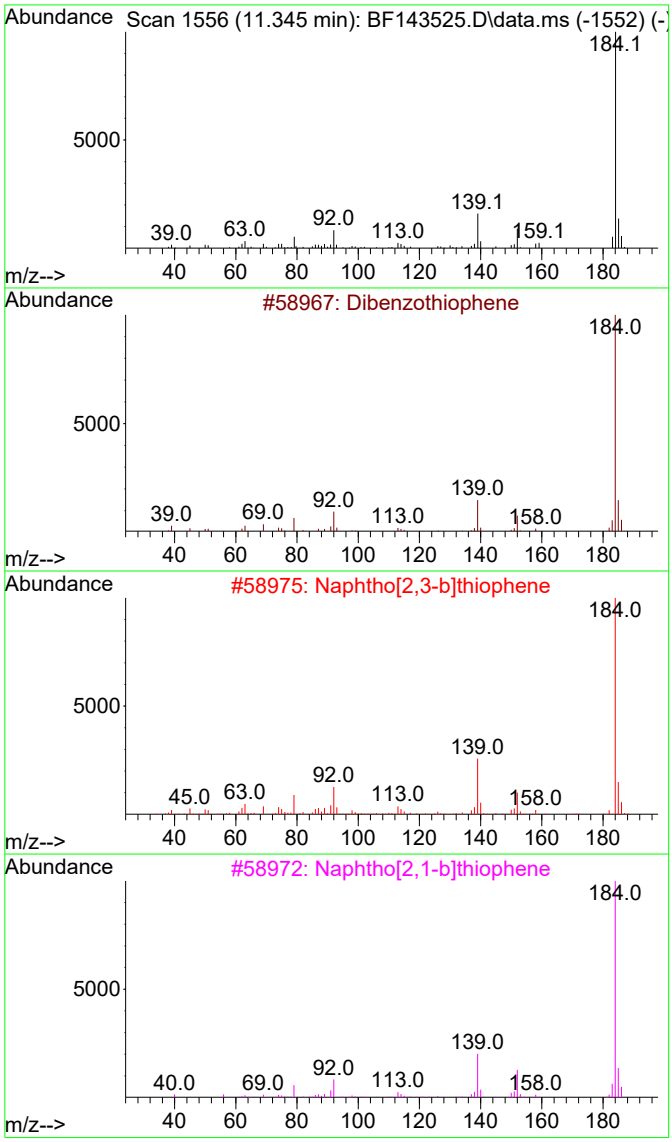
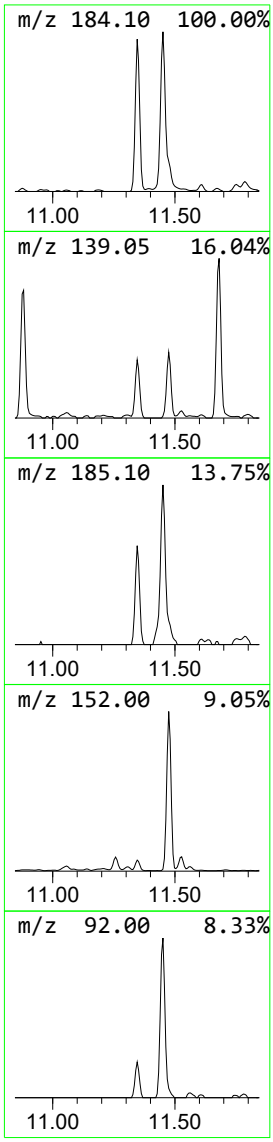
Peak Number 8 Dibenzothiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.345	3.09 ng	122567	Phenanthrene-d10	11.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	96
2	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	94
3	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	94
4	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91
5	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	90





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Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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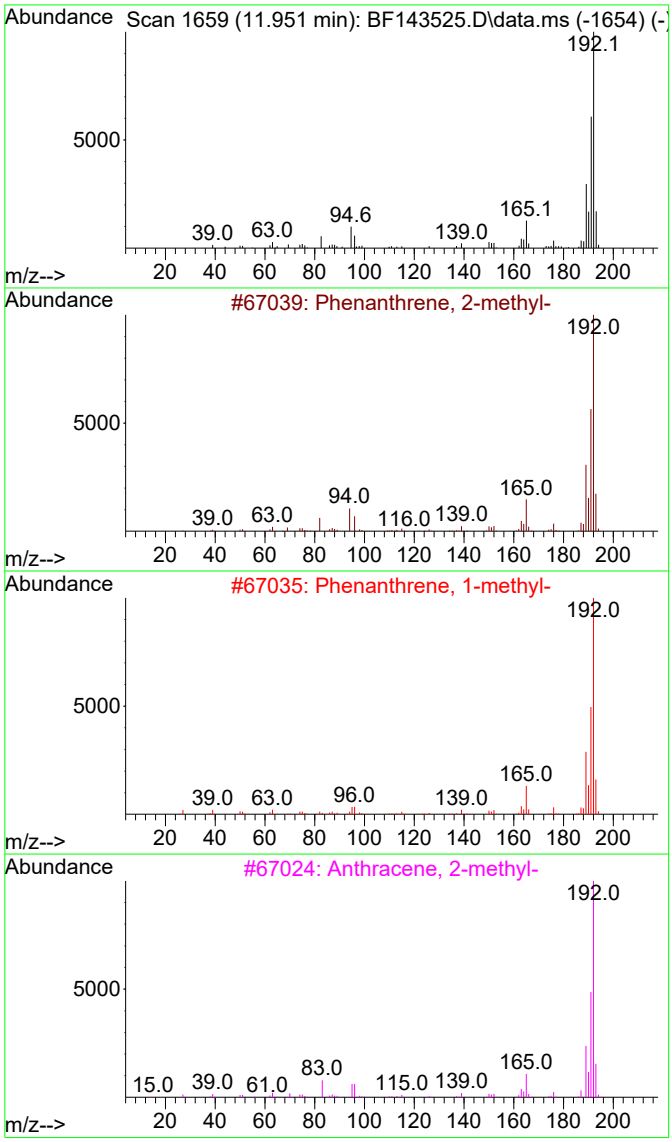
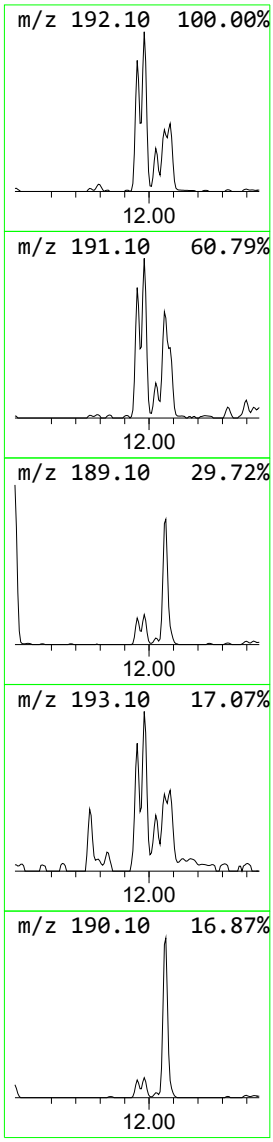
Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.951	2.56 ng	101599	Phenanthrene-d10	11.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93





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Acq On : 21 Aug 2025 12:46
Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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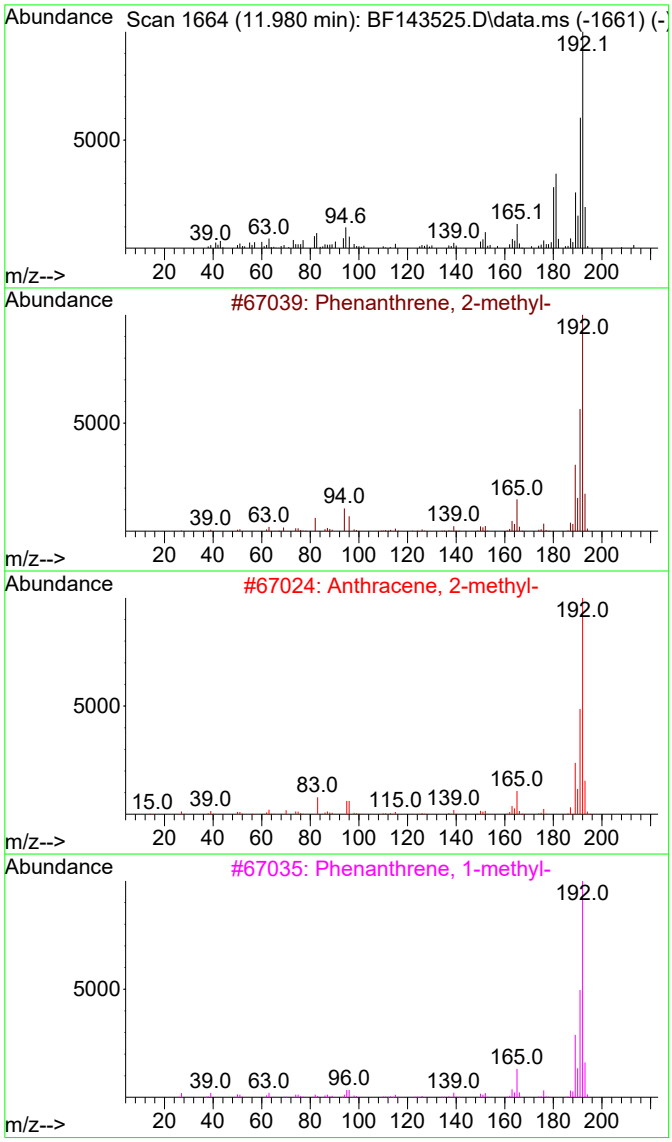
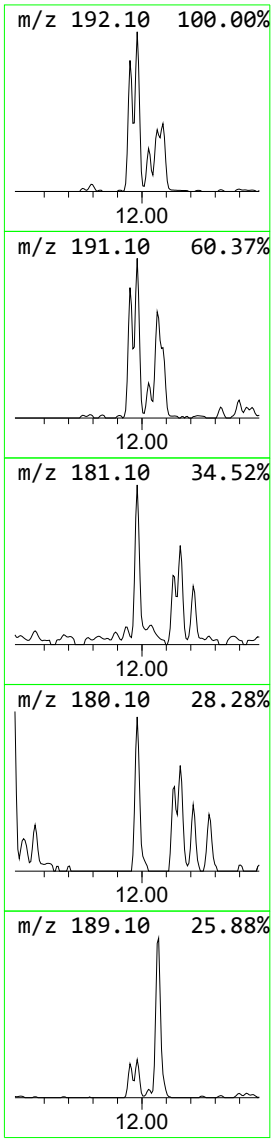
Peak Number 10 Anthracene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.980	4.14 ng	164367	Phenanthrene-d10	11.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	86





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Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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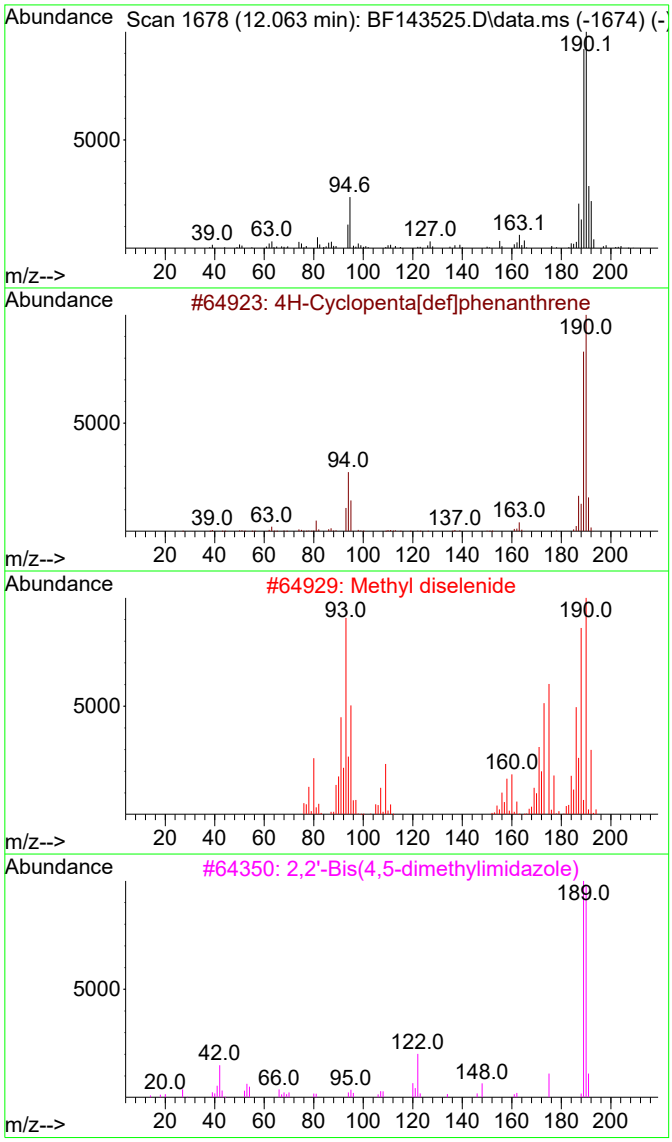
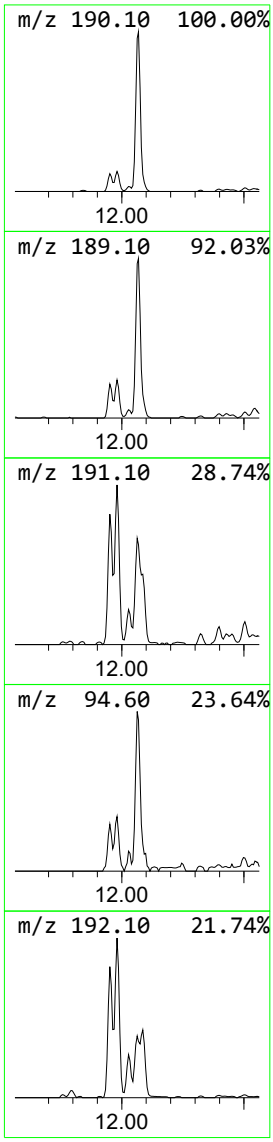
Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.063	6.31 ng	250377	Phenanthrene-d10	11.451

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	Methyl diselenide	190	C2H6Se2	007101-31-7	46
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
4	4-(4-Chlorophenyl)pyridine	189	C11H8ClN	005957-96-0	35
5	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25





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Acq On : 21 Aug 2025 12:46
Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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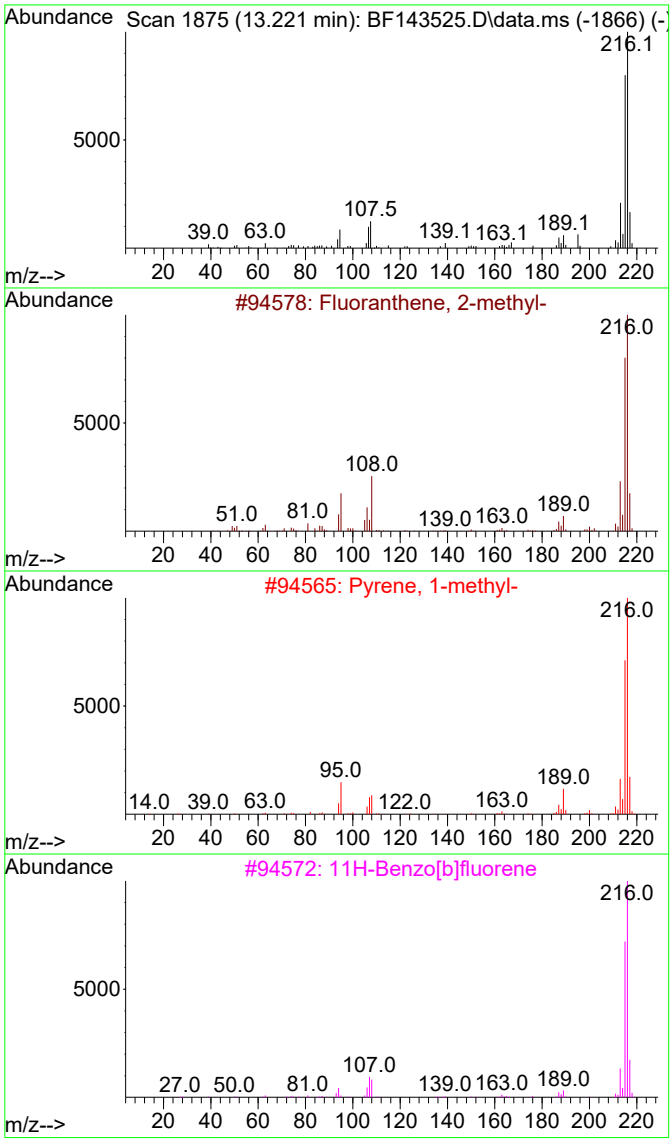
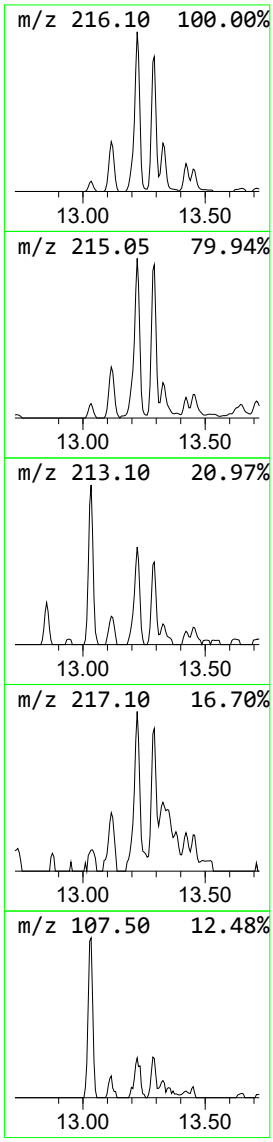
Peak Number 12 Fluoranthene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.222	2.70 ng	108678	Chrysene-d12	14.092

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	87





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Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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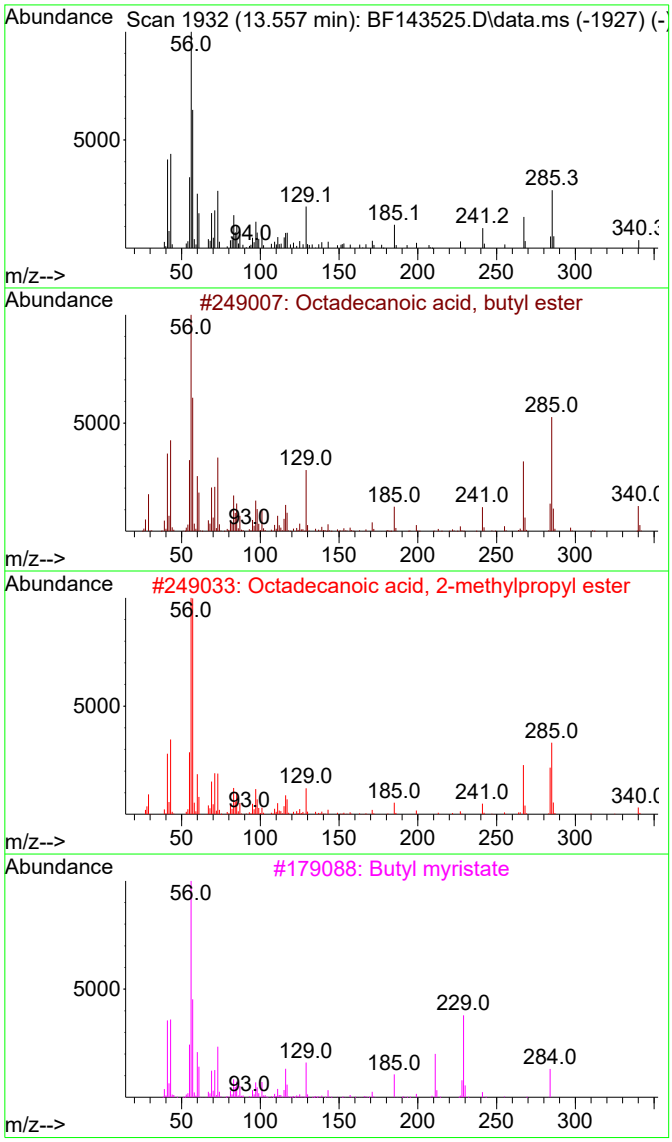
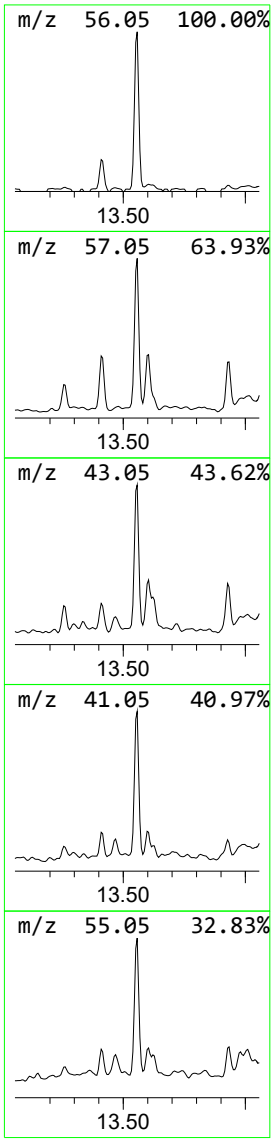
Peak Number 13 Octadecanoic acid, butyl ester Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.557	2.35 ng	94828	Chrysene-d12	14.092

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	99
2	Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	89
3	Butyl myristate	284	C18H36O2	000110-36-1	53
4	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	32
5	1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	30





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082125\
Data File : BF143525.D
Acq On : 21 Aug 2025 12:46
Operator : RC/JU
Sample : Q2819-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
11M-N

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	2.275	37.5	ng	1313710	1	6.928	700686	20.0
2-Pentanone, 4-...	5.157	6.4	ng	222905	1	6.928	700686	20.0
Naphthalene, 1,...	9.616	3.6	ng	188908	3	9.963	1058770	20.0
1,1'-Biphenyl, ...	10.592	2.1	ng	113894	3	9.963	1058770	20.0
Benzophenone	10.669	6.9	ng	362902	3	9.963	1058770	20.0
1(2H)-Acenaphth...	10.875	3.1	ng	121610	4	11.451	793426	20.0
9H-Fluorene, 1-...	11.057	2.6	ng	101668	4	11.451	793426	20.0
Dibenzothiophene	11.345	3.1	ng	122567	4	11.451	793426	20.0
Phenanthrene, 2...	11.951	2.6	ng	101599	4	11.451	793426	20.0
Anthracene, 2-m...	11.980	4.1	ng	164367	4	11.451	793426	20.0
4H-Cyclopenta[d...	12.063	6.3	ng	250377	4	11.451	793426	20.0
Fluoranthene, 2...	13.222	2.7	ng	108678	5	14.092	806377	20.0
Octadecanoic ac...	13.557	2.4	ng	94828	5	14.092	806377	20.0