LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

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_M\Methods\8270-BM041723.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM039564.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.301	210	215	222	rBV	112142	155442	2.18%	0.293%
2	4.913	313	319	329	rBV	249815	359265	5.03%	0.677%
3	5.384	393	399	424	rBV	2243618	3480483	48.76%	6.563%
4	5.795	464	469	479	rBV	41094	88680	1.24%	0.167%
5	7.001	664	674	695	rBV	2252611	3972047	55.65%	7.490%
6	7.472	749	754	764	rBV3	42797	93645	1.31%	0.177%
7	7.813	806	812	823	rBV	615750	1039488	14.56%	1.960%
8	8.360	899	905	916	rBV	50307	106722	1.50%	0.201%
9	8.989	1004	1012	1027	rBV	1427740	2533302	35.49%	4.777%
10	10.624	1279	1290	1296	rBV	811420	1456346	20.40%	2.746%
11	10.672	1296	1298	1306	rVB	67626	108372	1.52%	0.204%
12	12.295	1567	1574	1581	rBV2	69556	139512	1.95%	0.263%
13	12.513	1601	1611	1620	rVB	63074	122851	1.72%	0.232%
14	13.095	1703	1710	1729	rBV	3809344	5651347	79.17%	10.657%
15	13.795	1823	1829	1834	rBV	60590	106172	1.49%	0.200%
16	13.848	1834	1838	1847	rVB2	38088	72574	1.02%	0.137%
17	14.195	1892	1897	1909	rBV	140836	241753	3.39%	0.456%
18	14.471	1934	1944	1951	rBV	1231467	1886915	26.44%	3.558%
19	14.536	1951	1955	1965	rVB2	50717	104974	1.47%	0.198%
20	15.524	2118	2123	2135	rVB2	98740	179035	2.51%	0.338%
21	15.836	2169	2176	2189	rBV	561832	854339	11.97%	1.611%
22	15.965	2190	2198	2211	rBV	1845019	2823974	39.56%	5.325%
23	16.201	2231	2238	2249	rVB5	59777	167725	2.35%	0.316%
24	16.401	2267	2272	2283	rBV	94660	137818	1.93%	0.260%
25	17.224	2406	2412	2416	rBV	1416199	2035601	28.52%	3.838%
26	17.265	2416	2419	2424	rVV	838701	1194643	16.74%	2.253%
27	17.312	2424	2427	2430	rVV	115746	145930	2.04%	0.275%
28	17.359	2431	2435	2440	rVV	170974	256226	3.59%	0.483%
29	18.106	2557	2562	2563	rBV	201062	266979	3.74%	0.503%
30	18.153	2567	2570	2575	rVB	258687	375088	5.25%	0.707%
31	18.236	2581	2584	2589	rVB	65352	83240	1.17%	0.157%
32	18.300	2589	2595	2598	rVV2	223730	415539	5.82%	0.784%
33	18.336	2599	2601	2610	rVB	145547	187524	2.63%	0.354%
34	18.412	2611	2614	2619	rBV2	53674	77322	1.08%	0.146%
35	18.606	2643	2647	2651	rBV	97605	136851	1.92%	0.258%
36	19.024	2714	2718	2724	rBV2	134156	250612	3.51%	0.473%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

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_M\Methods\8270-BM041723.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.289	2758	2763	2769	rBV	1198718	1594011	22.33%	3.006%
38	19.653	2816	2825	2834	rBV	1278587	2108508	29.54%	3.976%
39	19.853	2854	2859	2869	rVB	5622604	7137908	100.00%	13.460%
40	20.147	2907	2909	2915	rVB2	280799	396664	5.56%	0.748%
41	20.247	2923	2926	2930	rVB	162975	189731	2.66%	0.358%
42	20.306	2933	2936	2940	rVB	177904	220669	3.09%	0.416%
43	20.447	2957	2960	2964	rBV	130284	166388	2.33%	0.314%
44	20.494	2965	2968	2973	rVV	129615	177183	2.48%	0.334%
45	21.141	3073	3078	3085	rVB4	285017	577969	8.10%	1.090%
46	21.324	3107	3109	3113	rVB	214573	213392	2.99%	0.402%
47	21.418	3118	3125	3129	rVV2	1899147	3492166	48.92%	6.585%
48	21.453	3129	3131	3140	rVB	752449	922341	12.92%	1.739%
49	23.059	3398	3404	3409	rBV2	526597	1321298	18.51%	2.492%
50	23.571	3488	3491	3498	rVB	374702	583410	8.17%	1.100%
51	23.777	3521	3526	3532	rBV2	979976	1746492	24.47%	3.293%
52	24.353	3620	3624	3635	rVB2	415137	874697	12.25%	1.649%

Sum of corrected areas: 53031163

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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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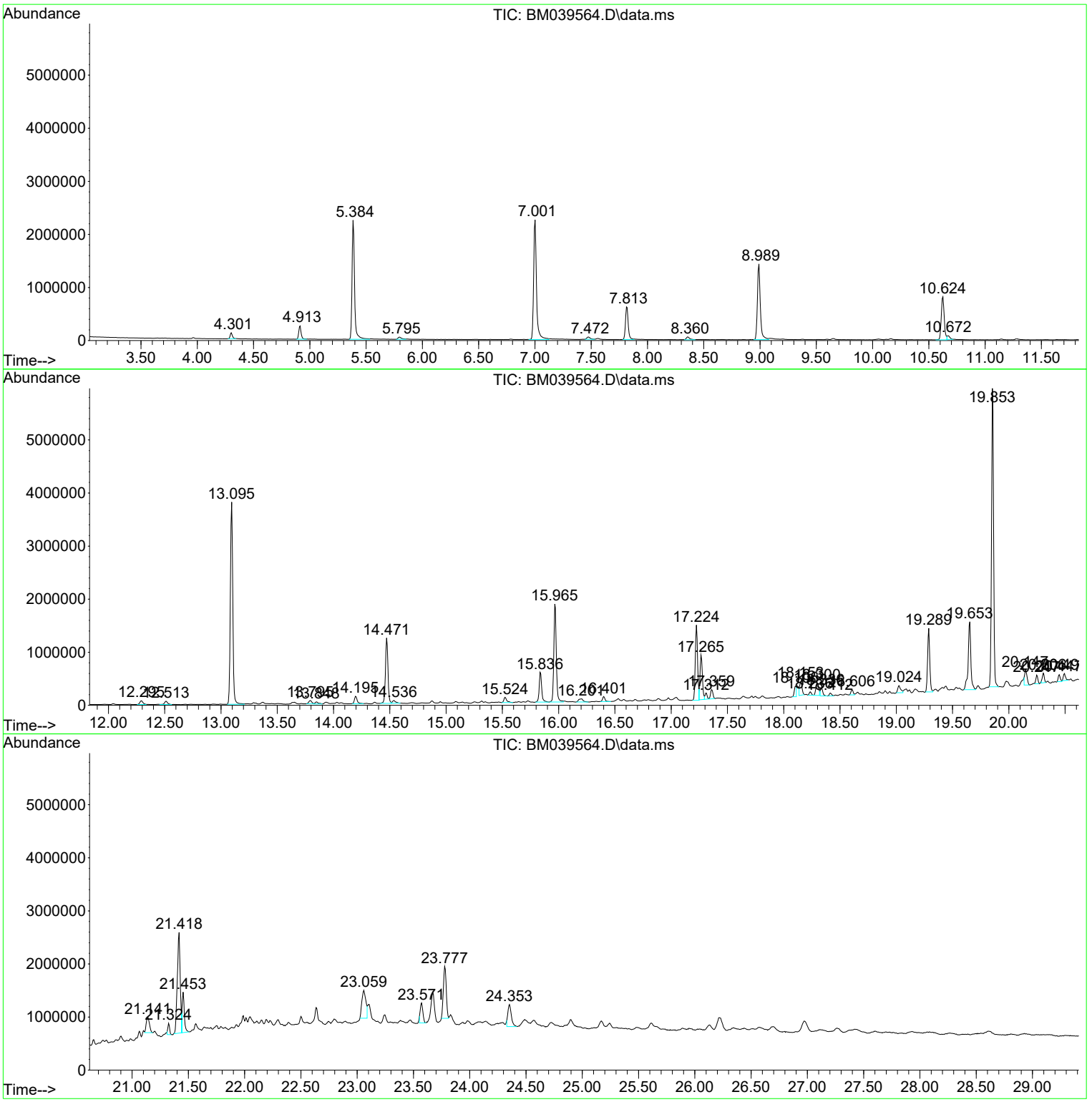
ClientSampleId :
NATIONAL-GRID-COMP

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

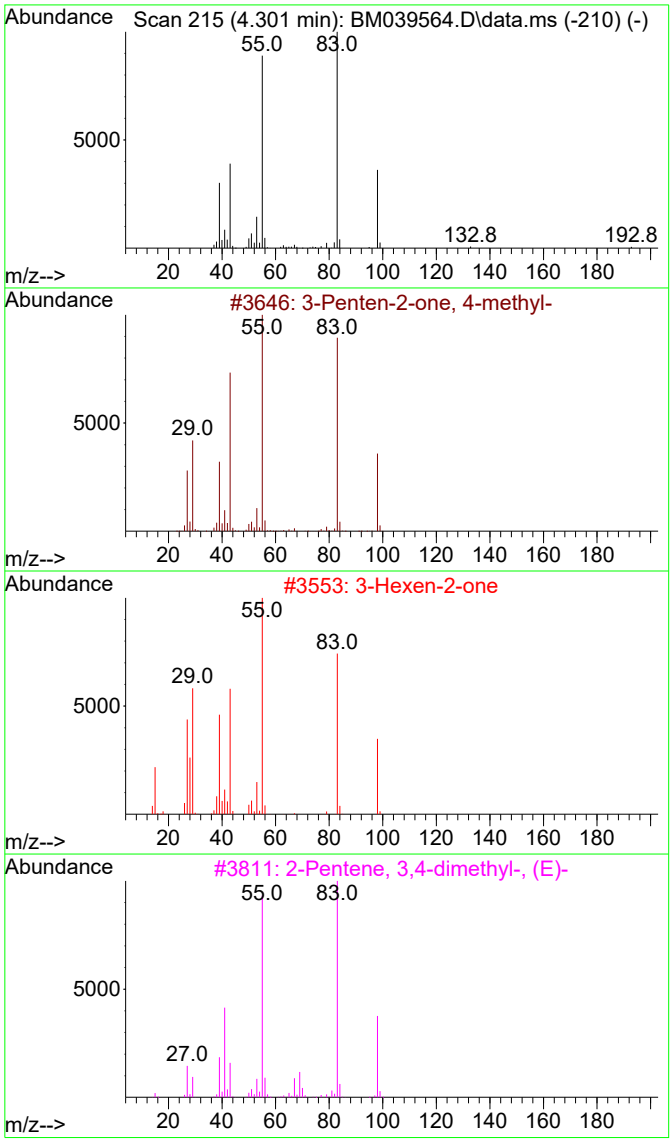
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

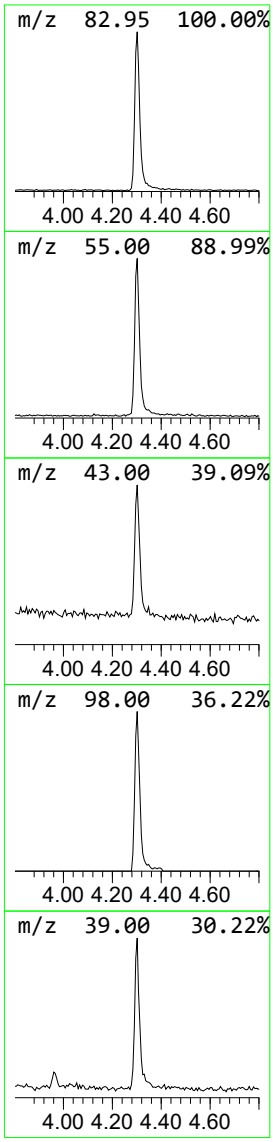
****

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.301	2.99 ng	155442	1,4-Dichlorobenzene-d4	7.813

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2	3-Hexen-2-one	98	C6H10O	000763-93-9	90
3	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	80
4	2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80
5	2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	72





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Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

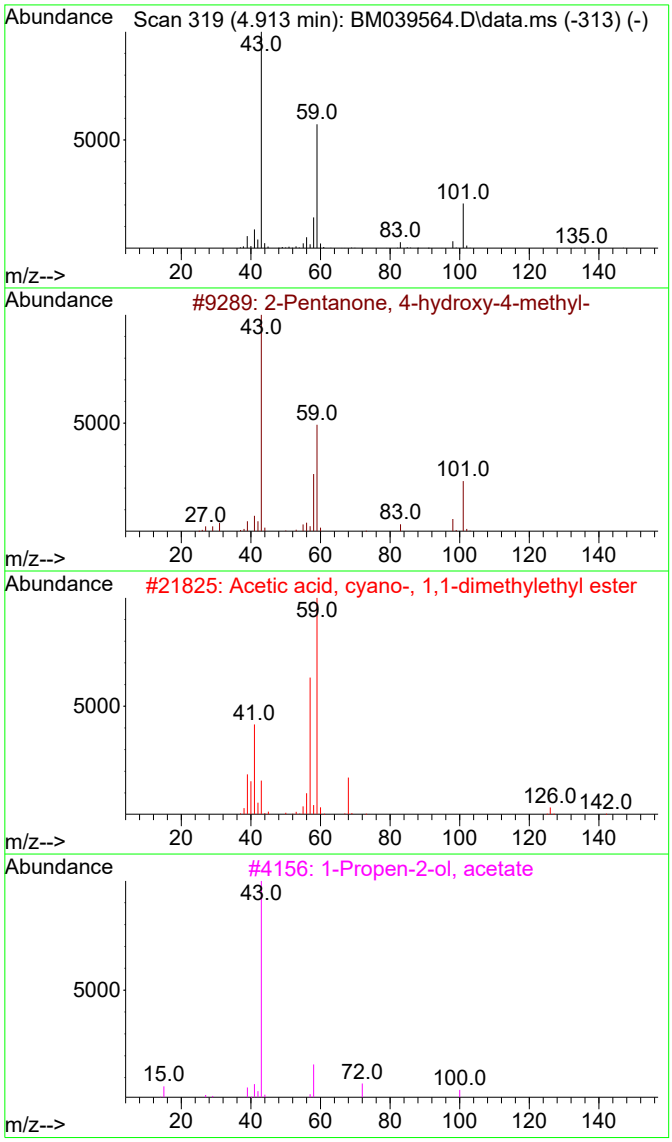
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

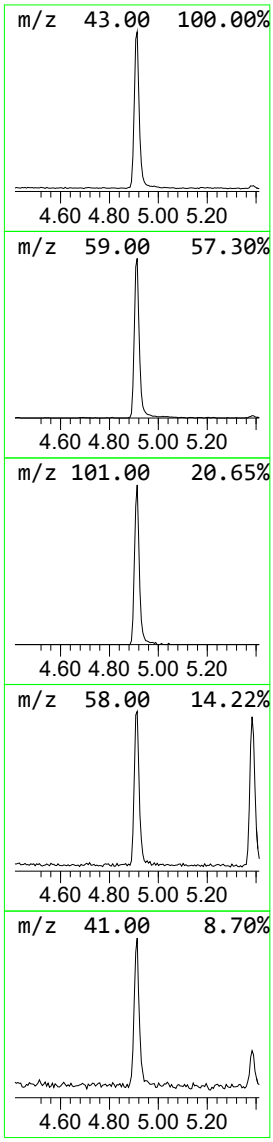
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.913	6.91 ng	359265	1,4-Dichlorobenzene-d4	7.813

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

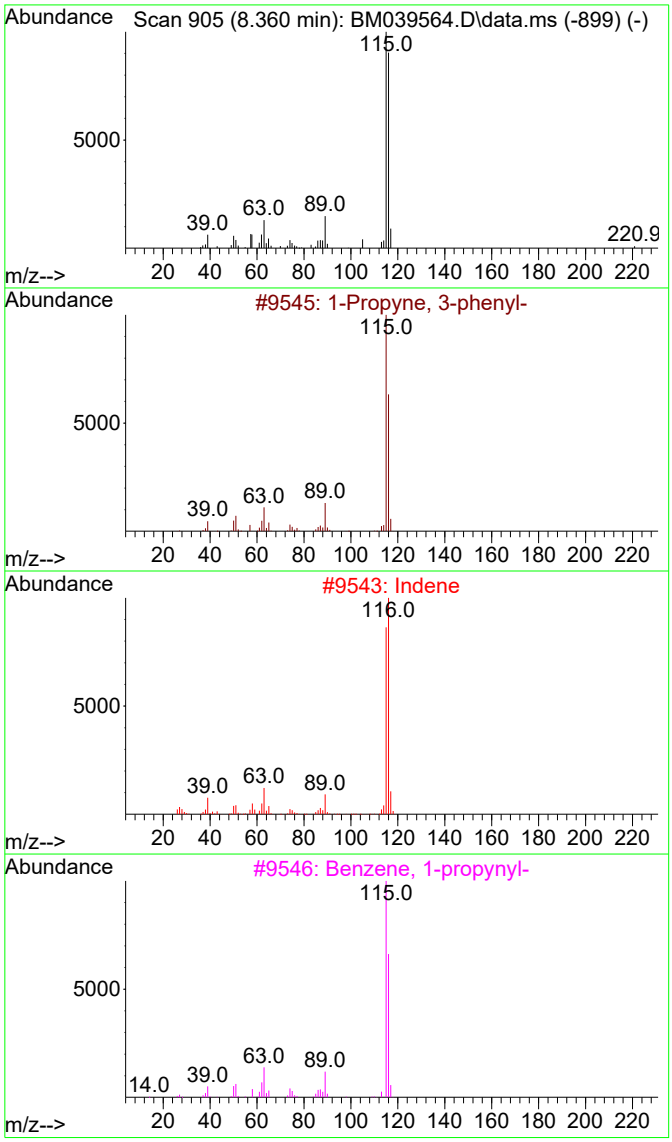
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

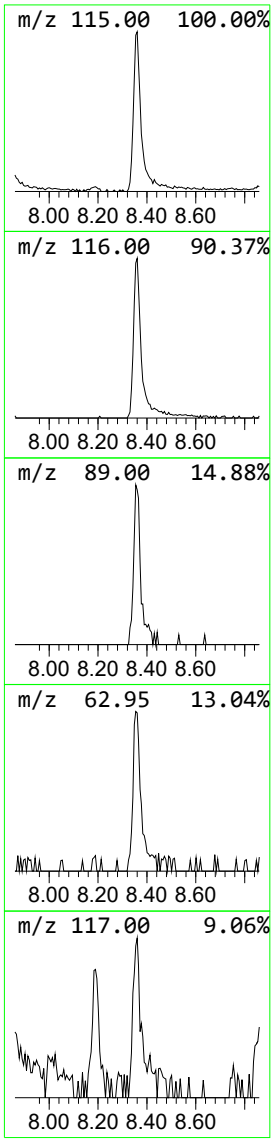
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Propyne, 3-phenyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.360	2.05 ng	106722	1,4-Dichlorobenzene-d4	7.813

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	97
2	Indene	116	C9H8	000095-13-6	94
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
4	3-Methylphenylacetylene	116	C9H8	000766-82-5	91
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91





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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

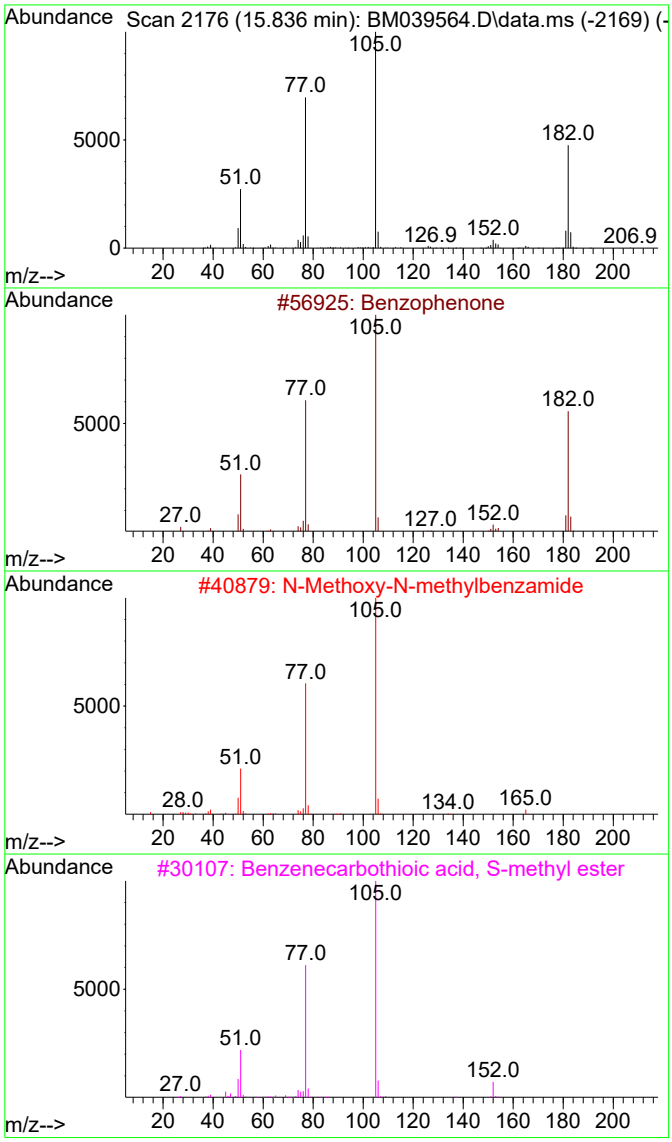
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

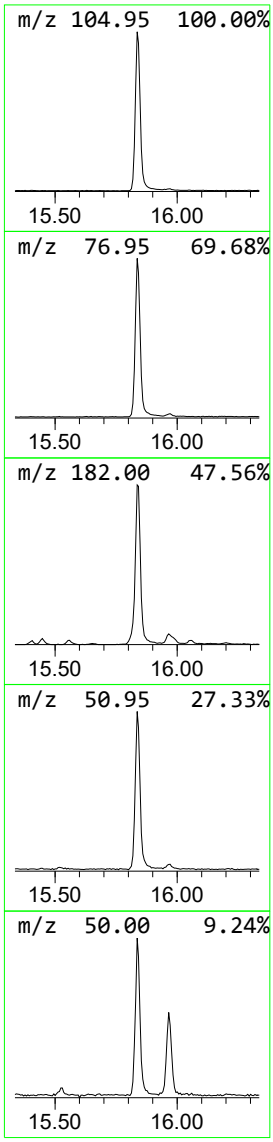
****

Peak Number 4 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.836	9.06 ng	854339	Acenaphthene-d10	14.471

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	58
3	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	52
4	2-Phenyl-4-ethylidene-2-oxazolin...	187	C11H9NO2	037791-41-6	50
5	acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	50





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ClientSampleId :
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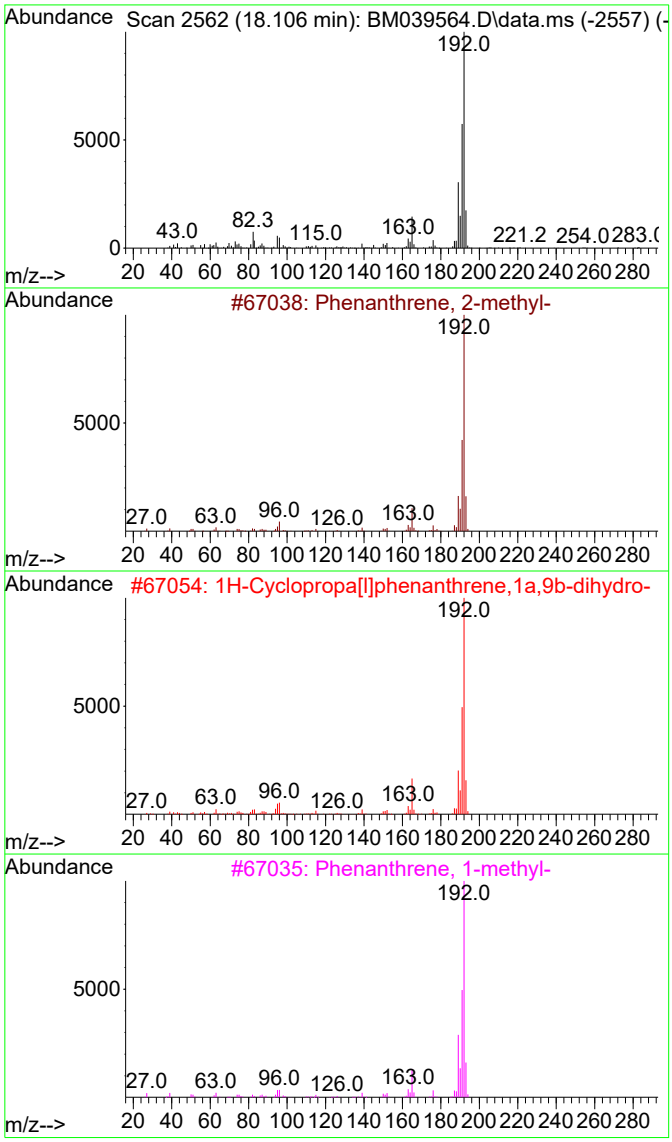
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

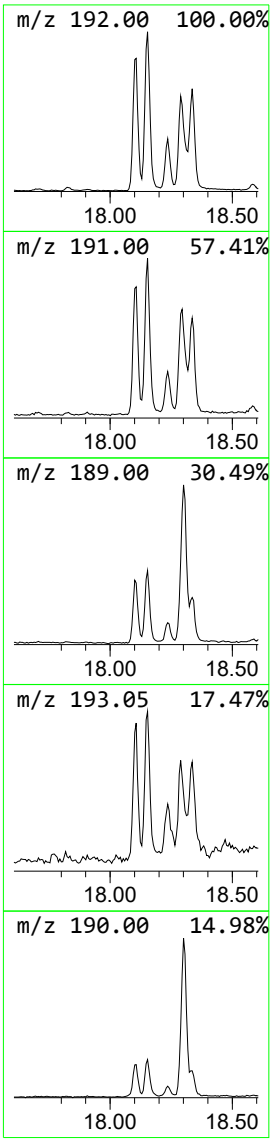
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.106	2.62 ng	266979	Phenanthrene-d10	17.224

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	97
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94





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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

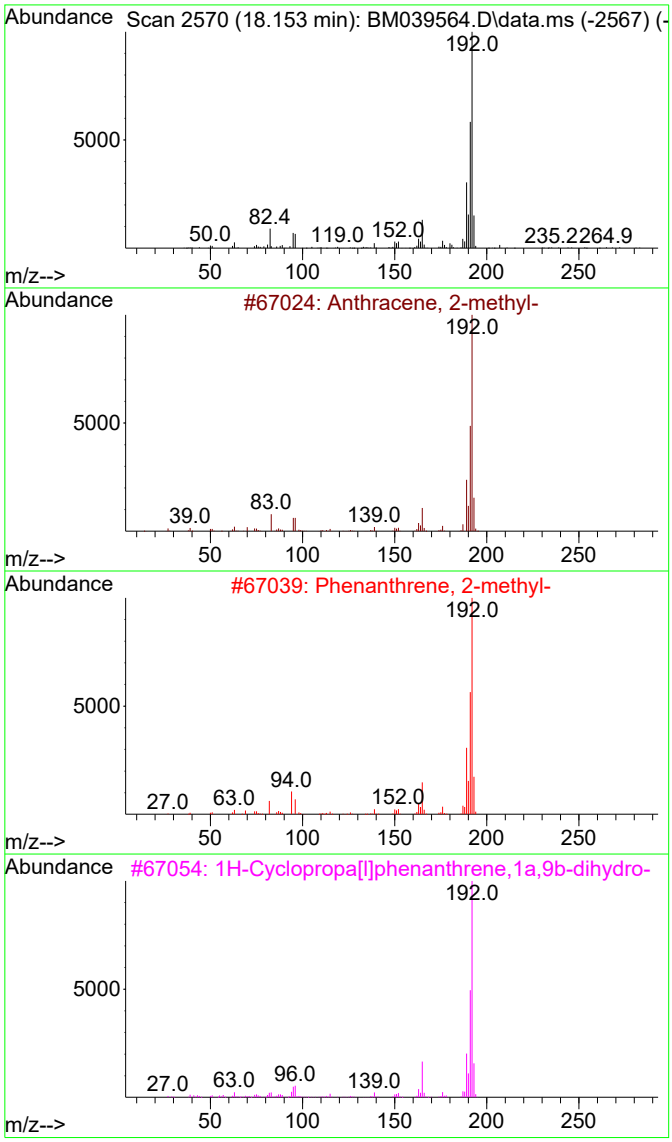
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

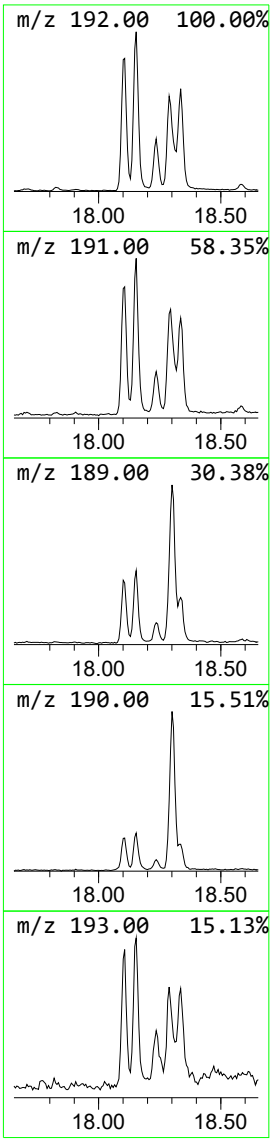
****

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.153	3.69 ng	375088	Phenanthrene-d10	17.224

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
4	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	94
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93





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Data File : BM039564.D
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Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

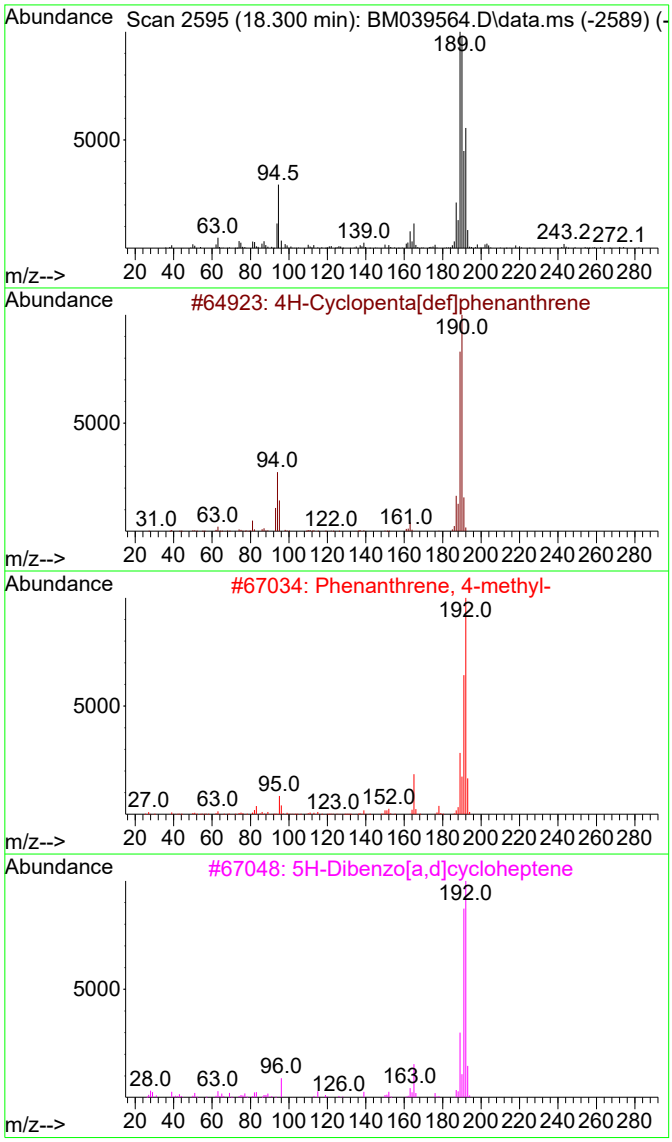
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

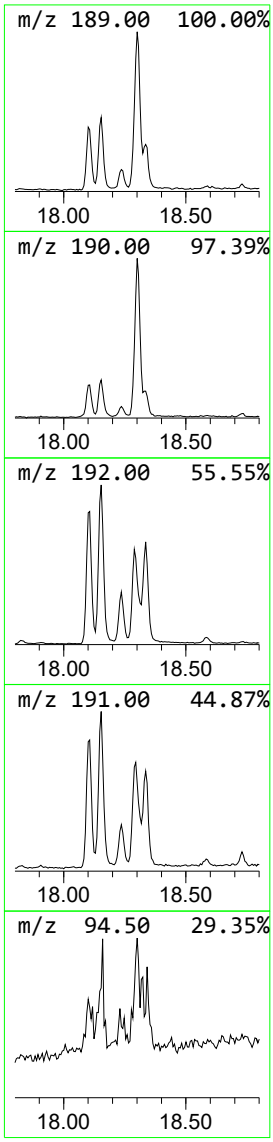
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.301	4.08 ng	415539	Phenanthrene-d10	17.224

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	35
3	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	18
4	Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	18
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	15





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

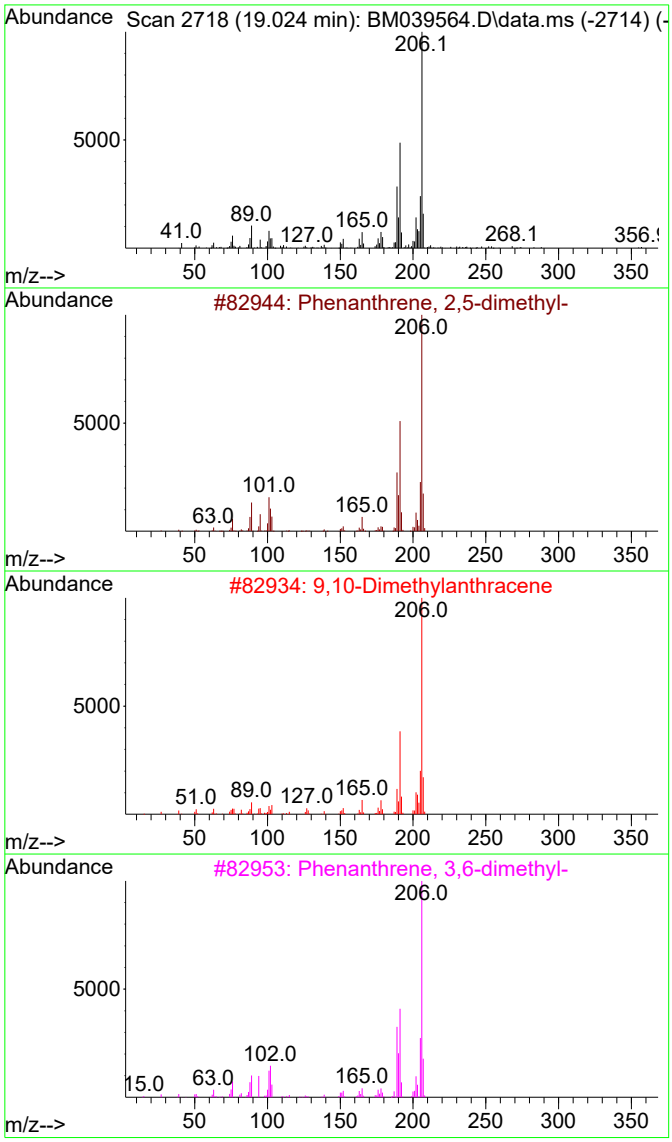
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

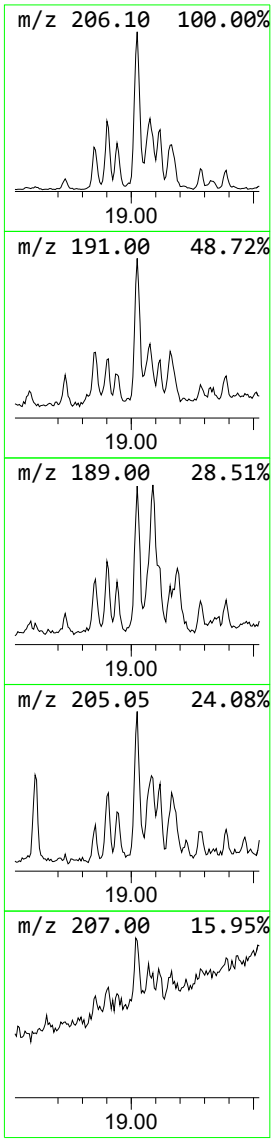
****

Peak Number 8 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.024	2.46 ng	250612	Phenanthrene-d10	17.224

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
3	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
4	di-p-Tolylacetylene	206	C16H14	002789-88-0	87
5	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86





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Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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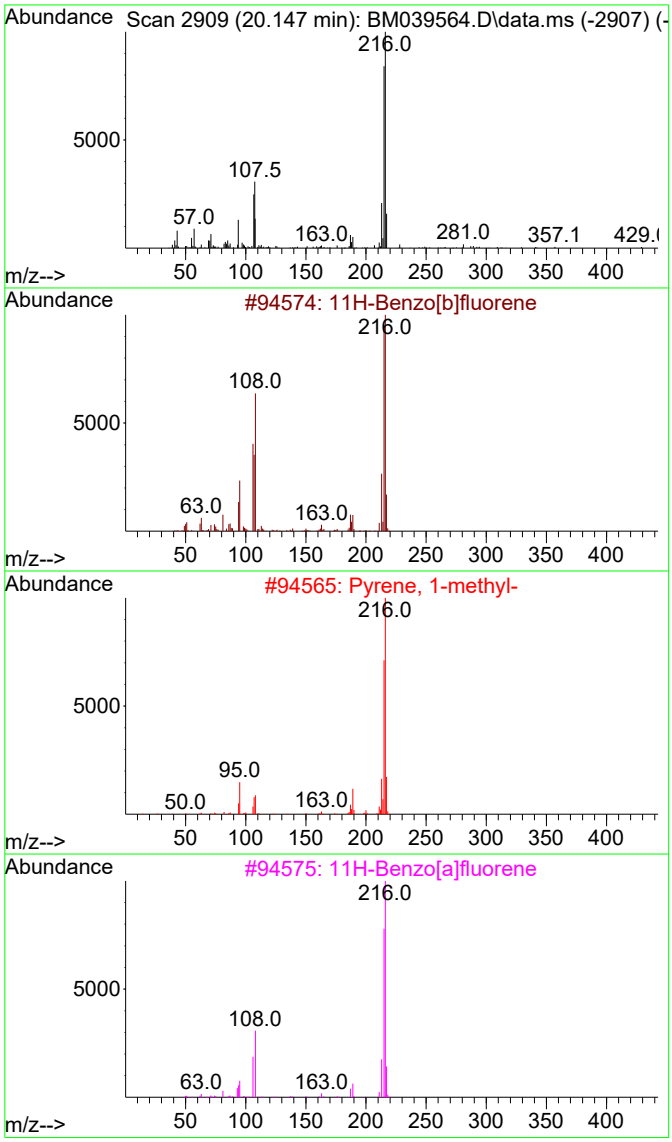
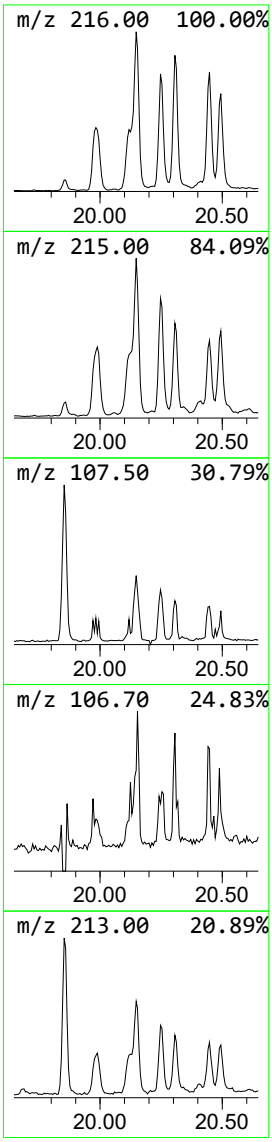
Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.147	2.27 ng	396664	Chrysene-d12	21.418

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	76
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	74
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	68
4	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	59
5	Allopregnane-3.alpha.,20.alpha.-...	320	C21H36O2	000566-58-5	50





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Sample : 02416-01
Misc :
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Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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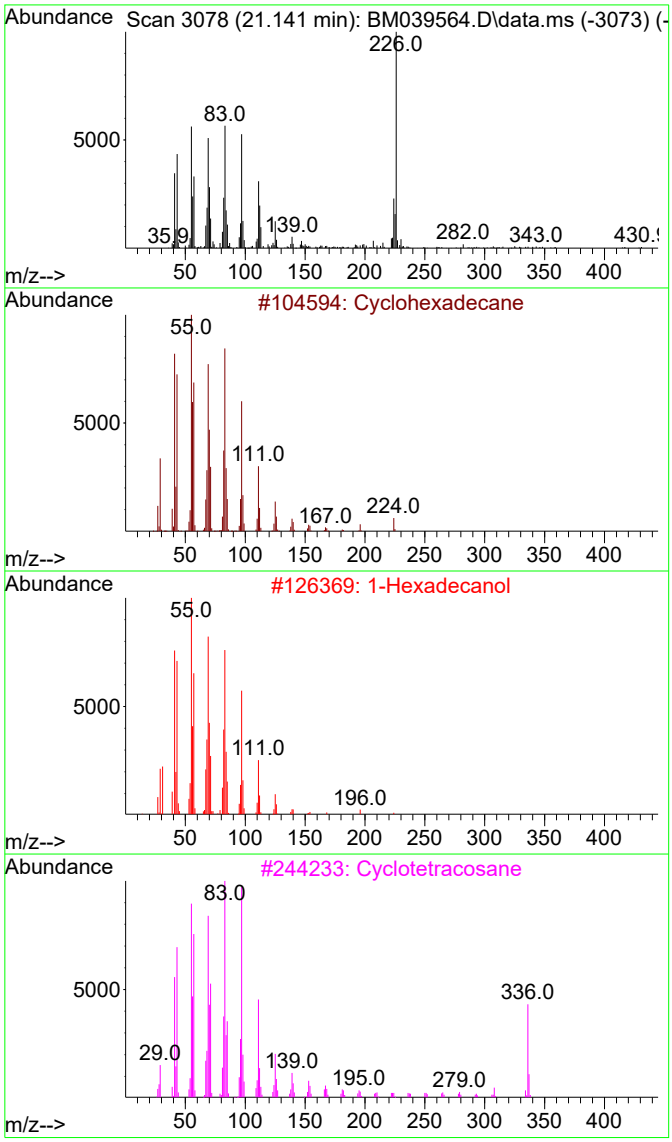
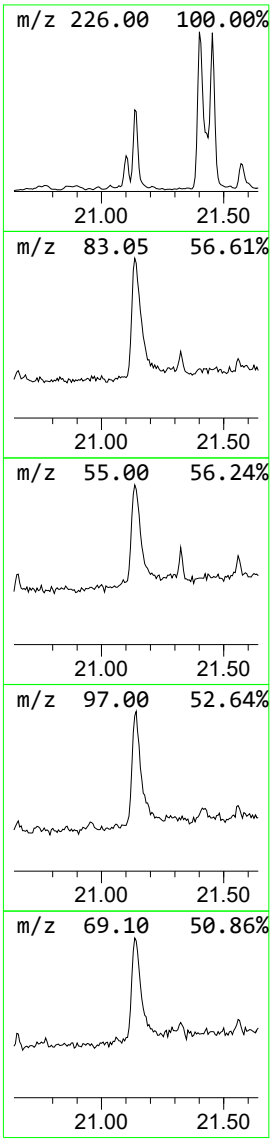
Peak Number 10 Cyclohexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.142	3.31 ng	577969	Chrysene-d12	21.418

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	96
2	1-Hexadecanol	242	C16H34O	036653-82-4	46
3	Cyclotetracosane	336	C24H48	000297-03-0	46
4	Behenic alcohol	326	C22H46O	000661-19-8	38
5	7-Hexadecene, (Z)-	224	C16H32	035507-09-6	35





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Data File : BM039564.D
Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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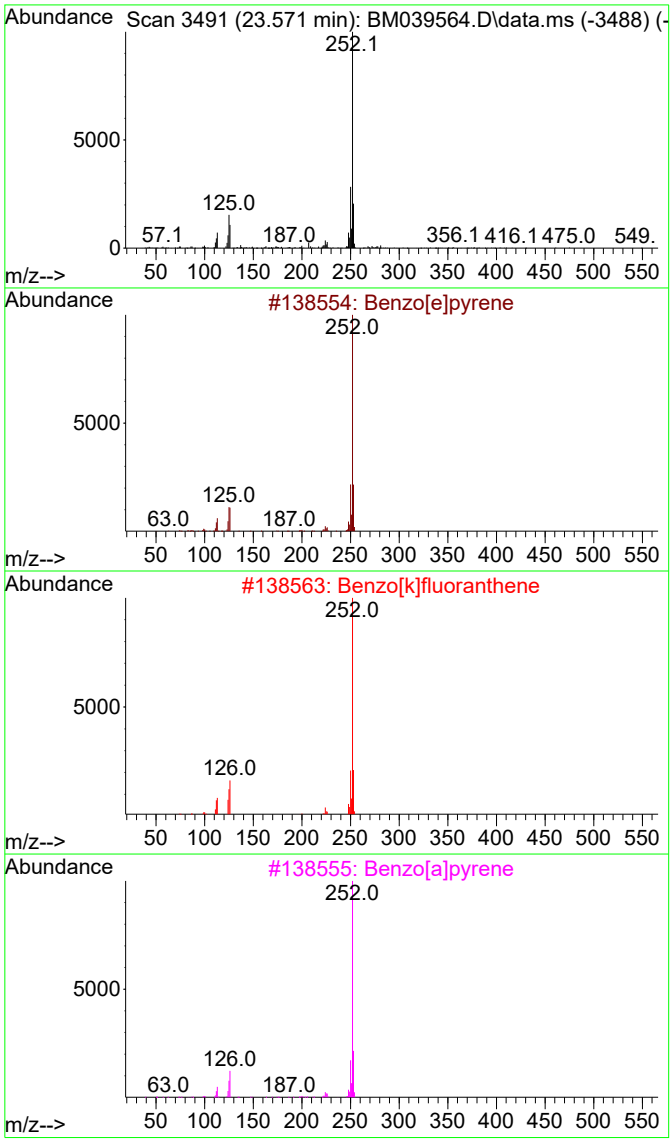
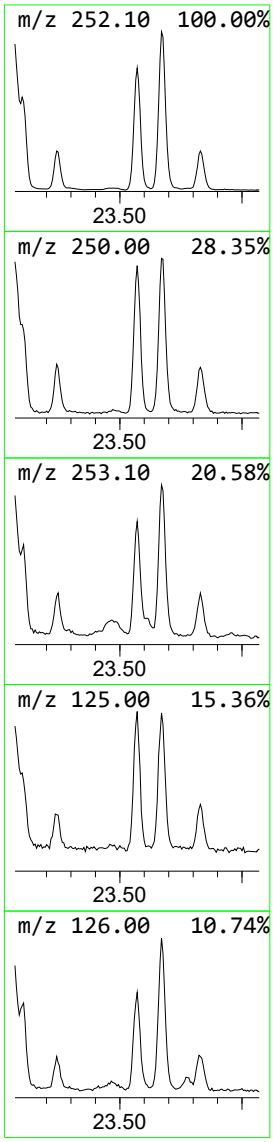
Peak Number 11 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.571	6.68 ng	583410	Perylene-d12	23.777

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3	Benzo[a]pyrene	252	C20H12	000050-32-8	96
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
5	Perylene	252	C20H12	000198-55-0	90





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Acq On : 21 Apr 2023 22:52
Operator : CG/JU
Sample : 02416-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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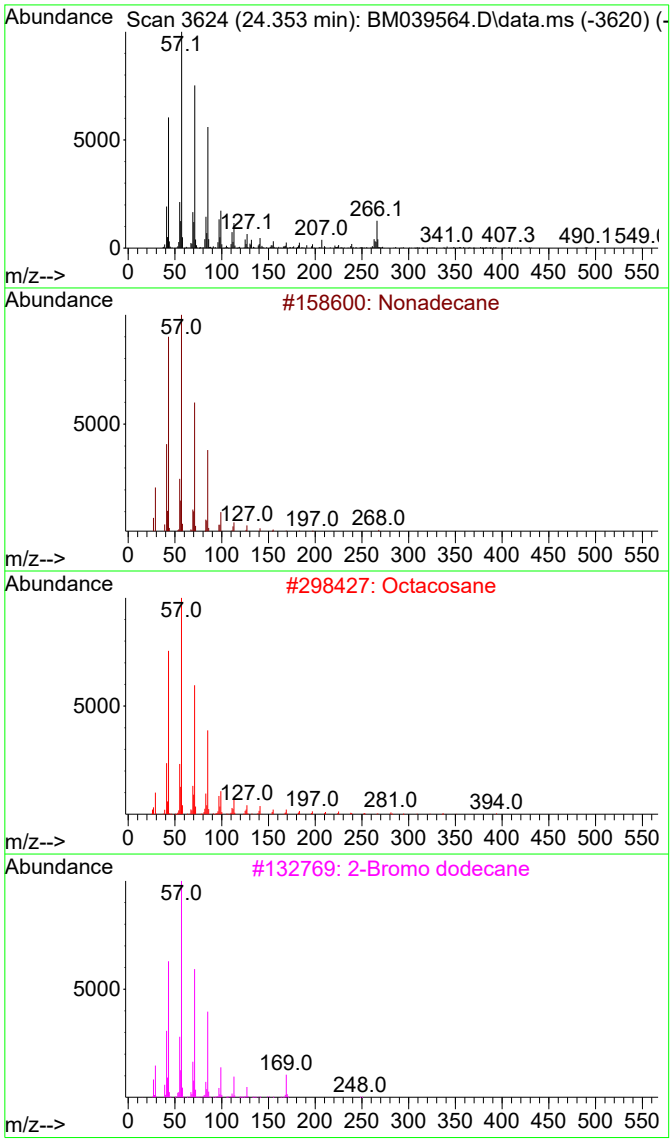
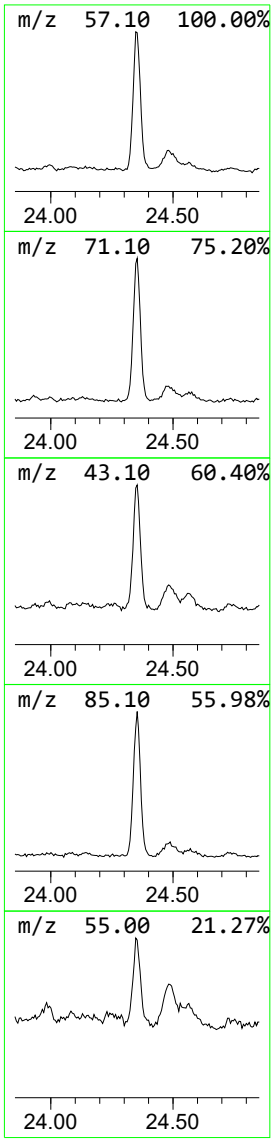
Peak Number 12 Nonadecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.353	10.02 ng	874697	Perylene-d12	23.777

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	96
2	Octacosane	394	C28H58	000630-02-4	90
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	89
4	Docosane, 11-butyl-	366	C26H54	013475-76-8	89
5	Heneicosane	296	C21H44	000629-94-7	87





Tentatively Identified Compound (LSC) summary

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Data File : BM039564.D
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Operator : CG/JU
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
NATIONAL-GRID-COMP

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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
3-Penten-2-one,...	4.301	3.0	ng	155442	1	7.813	1039490	20.0
2-Pentanone, 4-...	4.913	6.9	ng	359265	1	7.813	1039490	20.0
1-Propyne, 3-ph...	8.360	2.0	ng	106722	1	7.813	1039490	20.0
Benzophenone	15.836	9.1	ng	854339	3	14.471	1886920	20.0
Phenanthrene, 2...	18.106	2.6	ng	266979	4	17.224	2035600	20.0
Anthracene, 2-m...	18.153	3.7	ng	375088	4	17.224	2035600	20.0
4H-Cyclopenta[d...	18.301	4.1	ng	415539	4	17.224	2035600	20.0
Phenanthrene, 2...	19.024	2.5	ng	250612	4	17.224	2035600	20.0
11H-Benzo[b]flu...	20.147	2.3	ng	396664	5	21.418	3492170	20.0
Cyclohexadecane	21.142	3.3	ng	577969	5	21.418	3492170	20.0
Benzo[e]pyrene	23.571	6.7	ng	583410	6	23.777	1746490	20.0
Nonadecane	24.353	10.0	ng	874697	6	23.777	1746490	20.0