

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090623\
 Data File : BP016966.D
 Acq On : 06 Sep 2023 21:54
 Operator : MA/JU
 Sample : 04148-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBK25

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP090523.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016966.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.329	3	7	21	rVB	139808	223237	1.34%	0.143%
2	3.634	55	59	68	rVB2	34839	59705	0.36%	0.038%
3	3.770	78	82	99	rBV	444451	763248	4.57%	0.488%
4	3.964	110	115	122	rVB	58914	94952	0.57%	0.061%
5	4.081	130	135	140	rBV	37766	51829	0.31%	0.033%
6	4.470	197	201	205	rBV5	16451	26537	0.16%	0.017%
7	4.605	218	224	240	rVB	2964441	4543200	27.23%	2.904%
8	5.023	291	295	299	rBV	19708	26632	0.16%	0.017%
9	5.946	448	452	459	rVB2	26906	41563	0.25%	0.027%
10	6.381	520	526	530	rBV3	13423	20853	0.12%	0.013%
11	6.428	530	534	542	rVB2	19102	33282	0.20%	0.021%
12	6.928	613	619	623	rBV2	20201	33771	0.20%	0.022%
13	7.128	647	653	665	rBV	584991	918473	5.50%	0.587%
14	7.322	678	686	698	rBV	2331930	3615117	21.66%	2.311%
15	7.499	709	716	727	rBV	2109527	3363591	20.16%	2.150%
16	7.981	789	798	807	rBV	1955588	3025923	18.13%	1.934%
17	8.328	852	857	863	rBV8	12913	21759	0.13%	0.014%
18	8.687	910	918	929	rBV	1487576	2413626	14.46%	1.543%
19	8.899	948	954	964	rVB	84714	139396	0.84%	0.089%
20	9.181	994	1002	1007	rBV	2779048	4510621	27.03%	2.883%
21	9.228	1007	1010	1014	rVV	99206	159745	0.96%	0.102%
22	9.269	1014	1017	1027	rVB	53735	111957	0.67%	0.072%
23	9.393	1033	1038	1046	rVB2	19587	36036	0.22%	0.023%
24	9.893	1116	1123	1139	rBV	2369972	3780906	22.66%	2.417%
25	10.128	1157	1163	1172	rBV8	18063	40571	0.24%	0.026%
26	10.416	1205	1212	1225	rBV	3183177	5303255	31.78%	3.390%
27	10.658	1247	1253	1263	rVB7	11847	22265	0.13%	0.014%
28	10.810	1270	1279	1287	rBV	2651975	4535815	27.18%	2.900%
29	10.969	1297	1306	1320	rBV	858982	1496765	8.97%	0.957%
30	11.081	1320	1325	1334	rVB5	16600	36243	0.22%	0.023%
31	11.181	1334	1342	1348	rBV4	19895	41401	0.25%	0.026%
32	11.575	1402	1409	1412	rBV2	23364	40709	0.24%	0.026%
33	11.628	1412	1418	1425	rVB	135282	236559	1.42%	0.151%
34	11.810	1441	1449	1459	rBV	1318480	2472038	14.81%	1.580%
35	11.969	1469	1476	1484	rVB10	15444	39485	0.24%	0.025%
36	12.281	1525	1529	1534	rBV3	9408	16834	0.10%	0.011%

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Filtering: 5

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Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP090523.MA.M
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37	12.387	1541	1547	1556	rVB2	58342	99727	0.60%	0.064%
38	12.681	1592	1597	1602	rBV3	13772	24502	0.15%	0.016%
39	12.769	1609	1612	1619	rBV8	10685	20116	0.12%	0.013%
40	12.957	1638	1644	1655	rVB9	8040	24864	0.15%	0.016%
41	13.210	1681	1687	1692	rBV7	9795	17608	0.11%	0.011%
42	13.487	1726	1734	1738	rVV	88073	145000	0.87%	0.093%
43	13.599	1746	1753	1762	rVB	12935	40273	0.24%	0.026%
44	13.740	1771	1777	1781	rBV2	17818	29756	0.18%	0.019%
45	13.940	1806	1811	1818	rBV2	40688	62994	0.38%	0.040%
46	14.051	1823	1830	1838	rBV	6591773	9198127	55.12%	5.880%
47	14.340	1871	1879	1889	rBV	7376095	10545691	63.20%	6.741%
48	14.563	1912	1917	1923	rBV	38210	69525	0.42%	0.044%
49	14.640	1923	1930	1938	rVB	4248154	6045753	36.23%	3.865%
50	14.763	1946	1951	1955	rBV7	13536	24434	0.15%	0.016%
51	14.851	1960	1966	1979	rBV	669428	999264	5.99%	0.639%
52	15.634	2091	2099	2106	rBV	9568789	13606383	81.54%	8.698%
53	15.693	2106	2109	2114	rVV	140778	201429	1.21%	0.129%
54	15.757	2114	2120	2127	rVV	3375735	4544547	27.23%	2.905%
55	15.869	2136	2139	2144	rVB	42817	56543	0.34%	0.036%
56	16.145	2182	2186	2194	rBV4	46563	75236	0.45%	0.048%
57	16.298	2207	2212	2219	rBV2	48023	82502	0.49%	0.053%
58	16.369	2220	2224	2228	rVB2	36013	51288	0.31%	0.033%
59	17.134	2351	2354	2358	rVB4	41205	53609	0.32%	0.034%
60	17.404	2394	2400	2409	rBV	5104595	7149074	42.84%	4.570%
61	17.510	2411	2418	2435	rVB	10384409	14490302	86.84%	9.263%
62	17.663	2442	2444	2448	rVB2	30154	36403	0.22%	0.023%
63	17.834	2471	2473	2478	rVB	35549	42487	0.25%	0.027%
64	18.028	2502	2506	2515	rBV4	75276	142933	0.86%	0.091%
65	18.245	2539	2543	2551	rBV	467995	683437	4.10%	0.437%
66	18.510	2585	2588	2596	rVB3	41148	55170	0.33%	0.035%
67	18.769	2630	2632	2642	rVB5	60628	92951	0.56%	0.059%
68	19.063	2679	2682	2689	rVB	59289	87039	0.52%	0.056%
69	19.316	2721	2725	2729	rVV	128510	158398	0.95%	0.101%
70	19.381	2733	2736	2743	rVB	148394	205378	1.23%	0.131%
71	19.481	2749	2753	2762	rBV2	409122	599814	3.59%	0.383%
72	19.622	2774	2777	2781	rVB	48330	54803	0.33%	0.035%
73	19.745	2792	2798	2804	rBV	13089606	16687101	100.00%	10.667%
74	19.834	2810	2813	2817	rVB	209065	252268	1.51%	0.161%
75	20.251	2879	2884	2891	rBV	742529	971845	5.82%	0.621%
76	21.504	3092	3097	3104	rBV	5263477	6725216	40.30%	4.299%

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Sampling : 1 Min Area: 0.1 % 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 >
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77	23.880	3492	3501	3511	rVB2	6480562	13726130	82.26%	8.775%
78	24.045	3522	3529	3540	rVB	2933385	5952171	35.67%	3.805%

Sum of corrected areas: 156429990

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Instrument :
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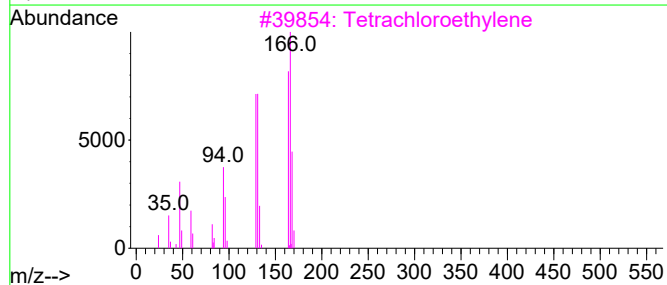
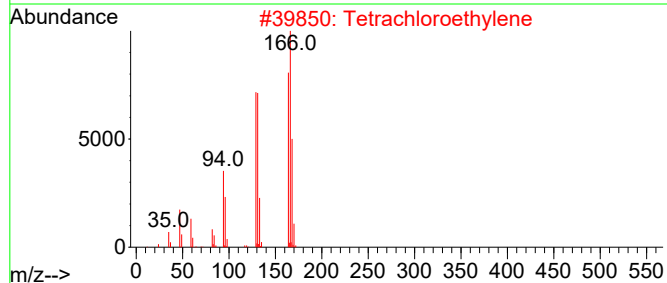
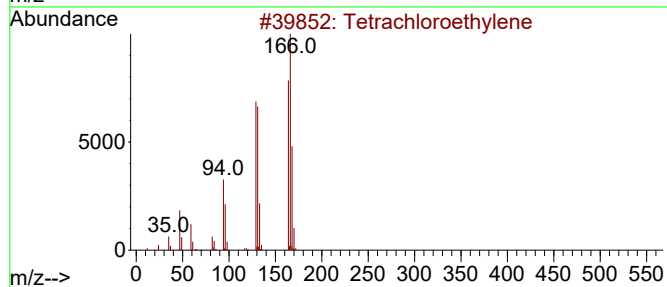
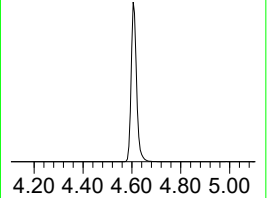
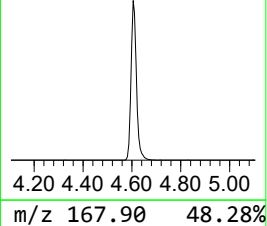
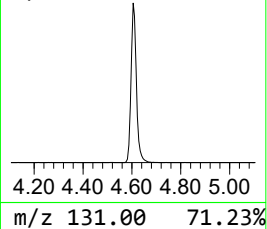
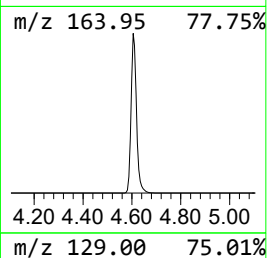
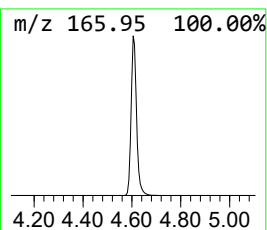
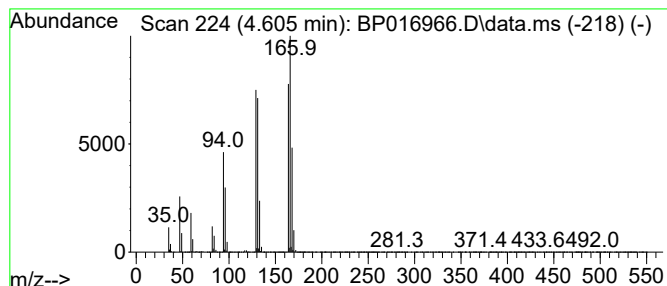
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.605	30.03 ng/ul	4543200	1,4-Dichlorobenzene-d4	7.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	94
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	35



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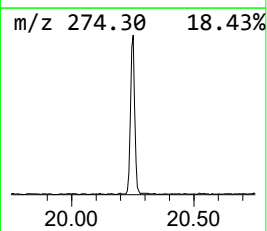
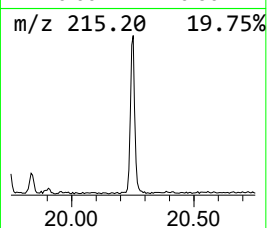
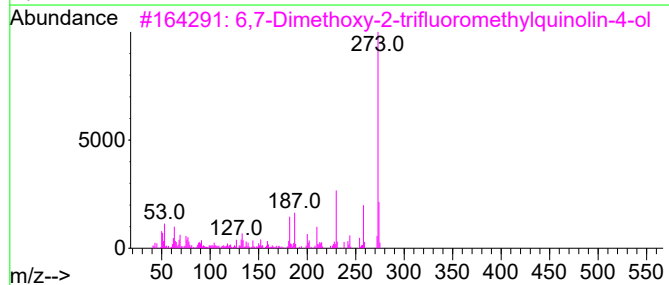
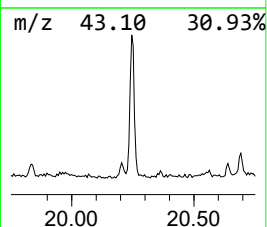
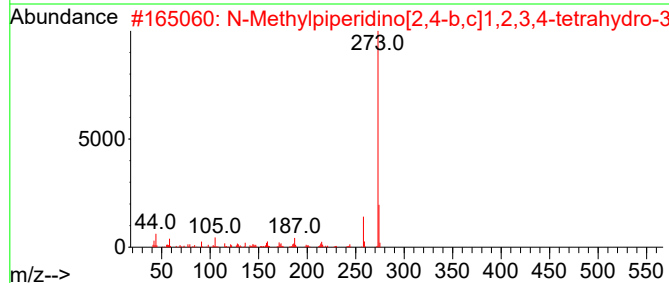
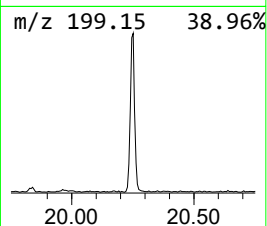
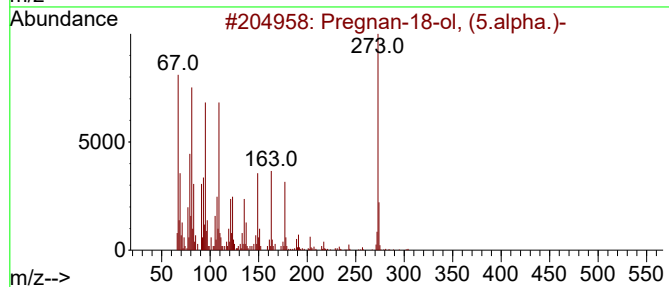
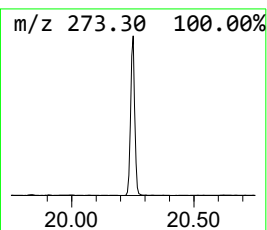
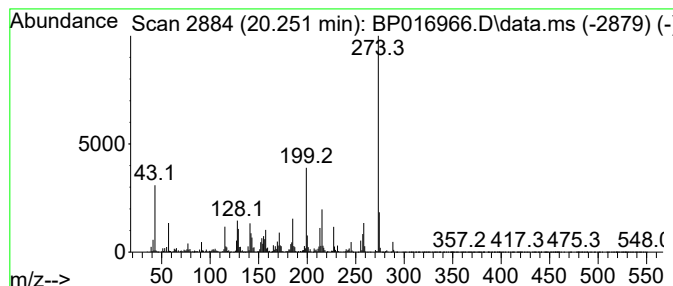
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.251	2.89 ng/ul	971845	Chrysene-d12	21.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pregnan-18-ol, (5.alpha.)-	304	C21H36O	056053-09-9	46
2			N-Methylpiperidino[2,4-b,c]1,2,3...	273	C18H27NO	1000128-54-7	45
3			6,7-Dimethoxy-2-trifluoromethylq...	273	C12H10F3NO3	041192-83-0	45
4			(R)-5-Methyl-6-(4-morpholinophen...	273	C15H19N3O2	1630760-58-5	43
5			3,3,6,6,9-Pentamethyl-3,4,5,6,7...	288	C18H24O3	019225-63-9	41



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.605	30.0	ng/ul	4543200	1	7.981	3025920	20.0
unknown-01	20.251	2.9	ng/ul	971845	5	21.504	6725220	20.0