

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011323\
 Data File : BN023643.D
 Acq On : 12 Jan 2023 14:03
 Operator : CG/JU
 Sample : N4239-04
 Misc : SFAM-MDL-S-02-4PPM
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-SOIL-QT3-2022-06

Quant Time: Jan 12 14:36:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010523.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 05 03:34:47 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 01/13/2023
 Supervised By :mohammad ahmed 01/13/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	182867	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.757	136	839389	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.598	164	535572	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	1178509	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1183129	20.000	ng/u1	-0.01
88) Perylene-d12	23.962	264	1196918	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	17072	4.017	ng/uL	0.00
4) Pyridine-d5	3.722	84	162099	12.714	ng/u1	0.00
7) Phenol-d5	7.105	99	288214	17.783	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	185861	18.694	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	230801	18.594	ng/u1	0.00
15) 4-Methylphenol-d8	8.657	113	232611	17.982	ng/u1	-0.01
21) Nitrobenzene-d5	9.116	128	114843	18.144	ng/u1	0.00
24) 2-Nitrophenol-d4	9.840	143	120075	17.920	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	226242	17.702	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.898	131	342364	17.030	ng/u1	-0.01
46) Dimethylphthalate-d6	14.004	166	698270	17.527	ng/u1	-0.01
49) Acenaphthylene-d8	14.292	160	794658	18.002	ng/u1	0.00
54) 4-Nitrophenol-d4	14.792	143	131313	16.190	ng/u1	-0.01
60) Fluorene-d10	15.586	176	635846	18.678	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.710	200	102166	14.371	ng/u1	0.00
73) Anthracene-d10	17.445	188	958852	18.182	ng/u1	0.00
81) Pyrene-d10	19.739	212	1185636	17.522	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.798	264	1083850	18.311	ng/u1	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.328	88	5851m	1.304	ng/uL	
5) Pyridine	3.752	79	48912	3.747	ng/u1	90
6) Benzaldehyde	7.075	77	35698	5.687	ng/u1	96
8) Phenol	7.128	94	68750	4.167	ng/u1	99
10) Bis(2-Chloroethyl)ether	7.363	93	56424	4.328	ng/u1	98
12) 2-Chlorophenol	7.499	128	55176	4.302	ng/u1	99
13) 2-Methylphenol	8.393	108	48546	3.910	ng/u1	93
14) 2,2'-oxybis(1-Chloropr...	8.481	45	78244	4.488	ng/u1	99
16) Acetophenone	8.775	105	86310	4.317	ng/u1	99
17) N-Nitroso-di-n-propyla...	8.757	70	41779	4.150	ng/u1	98
18) 4-Methylphenol	8.722	108	55593	4.066	ng/u1	99
19) Hexachloroethane	9.028	117	21584	4.207	ng/u1	99
22) Nitrobenzene	9.157	77	67081	4.235	ng/u1	98
23) Isophorone	9.681	82	107309	3.745	ng/u1	100
25) 2-Nitrophenol	9.869	139	30109	4.043	ng/u1	91
26) 2,4-Dimethylphenol	9.928	107	55602	3.679	ng/u1	97
27) Bis(2-Chloroethoxy)met...	10.169	93	76783	4.132	ng/u1	99
29) 2,4-Dichlorophenol	10.404	162	53054	4.149	ng/u1	99
30) Naphthalene	10.810	128	195118	4.324	ng/u1	99
32) 4-Chloroaniline	10.922	127	82300	4.088	ng/u1	99
33) Hexachlorobutadiene	11.093	225	33147	4.191	ng/u1	96
34) Caprolactam	11.710	113	16006	3.727	ng/u1	99
35) 4-Chloro-3-methylphenol	12.045	107	54317	3.895	ng/u1	99
36) 2-Methylnaphthalene	12.422	142	127753	4.247	ng/u1	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.640	142	118146	3.844	ng/u1	99
39) 1,2,4,5-Tetrachloroben...	12.787	216	66389	4.191	ng/u1	99
40) Hexachlorocyclopentadiene	12.769	237	35774	3.583	ng/u1	98
41) 2,4,6-Trichlorophenol	13.028	196	38261	3.664	ng/u1	97
42) 2,4,5-Trichlorophenol	13.098	196	43796	3.836	ng/u1	93
43) 1,1'-Biphenyl	13.428	154	171716	4.156	ng/u1	99
44) 2-Chloronaphthalene	13.469	162	136783	4.267	ng/u1	99
45) 2-Nitroaniline	13.681	65	30513	3.452	ng/u1	96
47) Dimethylphthalate	14.051	163	163835	4.092	ng/u1	99
48) 2,6-Dinitrotoluene	14.175	165	24323	3.214	ng/u1	91
50) Acenaphthylene	14.316	152	208334	4.245	ng/u1	99
51) 3-Nitroaniline	14.504	138	28427	3.748	ng/u1	94
52) Acenaphthene	14.657	153	146779	4.242	ng/u1	97
53) 2,4-Dinitrophenol	14.710	184	17656	3.569	ng/u1	98
55) 4-Nitrophenol	14.804	109	25820	3.942	ng/u1	93
56) Dibenzofuran	14.992	168	202964	4.214	ng/u1	93
57) 2,4-Dinitrotoluene	14.963	165	39131	3.542	ng/u1	98
58) 2,3,4,6-Tetrachlorophenol	15.222	232	36323	3.712	ng/u1	99
59) Diethylphthalate	15.416	149	154721	3.858	ng/u1	98
61) Fluorene	15.645	166	167791	4.302	ng/u1	97
62) 4-Chlorophenyl-phenyle...	15.639	204	80898	4.260	ng/u1	97
63) 4-Nitroaniline	15.663	138	29403	4.211	ng/u1	98
66) 4,6-Dinitro-2-methylph...	15.722	198	29570	4.215	ng/u1#	89
67) N-Nitrosodiphenylamine	15.851	169	136825	4.137	ng/u1	99
68) 4-Bromophenyl-phenylether	16.533	248	46862	3.967	ng/u1	96
69) Hexachlorobenzene	16.645	284	55652	4.020	ng/u1	94
70) Atrazine	16.804	200	47493	3.837	ng/u1	99
71) Pentachlorophenol	16.992	266	32667	3.698	ng/u1	94
72) Phenanthrene	17.386	178	267496	4.236	ng/u1	98
74) Anthracene	17.475	178	264571	4.210	ng/u1	99
75) 1,2,3,4-Tetrachloroben...	13.392	216	59871	3.773	ng/uL	92
76) Pentachlorobenzene	14.910	250	59240	3.727	ng/uL	97
77) Carbazole	17.751	167	235530	4.127	ng/u1	99
78) Di-n-butylphthalate	18.316	149	245848	3.629	ng/u1	99
80) Fluoranthene	19.404	202	309830	3.915	ng/u1	98
82) Pyrene	19.769	202	347541	4.158	ng/u1	99
83) Butylbenzylphthalate	20.669	149	112791	3.408	ng/u1	98
84) 3,3'-Dichlorobenzidine	21.451	252	85070	3.292	ng/u1	99
85) Benzo(a)anthracene	21.516	228	329794	4.115	ng/u1	97
86) Bis(2-ethylhexyl)phtha...	21.445	149	163676	3.406	ng/u1	97
87) Chrysene	21.569	228	320228	4.262	ng/u1	99
89) Di-n-octyl phthalate	22.374	149	250535	3.033	ng/u1	100
90) Benzo(b)fluoranthene	23.215	252	337767	4.430	ng/u1	98
91) Benzo(k)fluoranthene	23.263	252	315906	4.189	ng/u1	99
93) Benzo(a)pyrene	23.851	252	290621	4.395	ng/u1#	94
94) Indeno(1,2,3-cd)pyrene	26.480	276	346392	4.239	ng/u1	99
95) Dibenzo(a,h)anthracene	26.504	278	289496	4.284	ng/u1	99
96) Benzo(g,h,i)perylene	27.256	276	273134	4.197	ng/u1	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

