

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
 Data File : BP019402.D
 Acq On : 19 Jan 2024 15:15
 Operator : MA/JU
 Sample : PB158500BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLCS500

Quant Time: Jan 19 15:48:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 09 03:05:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	115293	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.563	136	463271	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.416	164	327186	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.181	188	771505	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	682573	20.000	ng/u1	0.00
88) Perylene-d12	23.645	264	694588	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	18591	6.098	ng/uL	-0.01
4) Pyridine-d5	3.622	84	263324	29.867	ng/u1	-0.02
7) Phenol-d5	6.952	99	294223	26.700	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.105	67	180524	27.389	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.299	132	211419	27.930	ng/u1	-0.02
15) 4-Methylphenol-d8	8.499	113	231792	26.514	ng/u1	-0.02
21) Nitrobenzene-d5	8.934	128	110074	28.242	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.652	143	127613	28.180	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.199	165	257118	28.103	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.716	131	274957	23.792	ng/u1	-0.01
46) Dimethylphthalate-d6	13.840	166	788396	27.275	ng/u1	0.00
49) Acenaphthylene-d8	14.110	160	845736	27.294	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.663	143	109435	24.654	ng/u1	0.00
60) Fluorene-d10	15.416	176	672060	28.398	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	157248	27.607	ng/u1	0.00
73) Anthracene-d10	17.281	188	1048149	27.240	ng/u1	0.00
81) Pyrene-d10	19.528	212	1259506	28.298	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.492	264	1063145	28.149	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.246	88	40937	12.962	ng/uL	98
5) Pyridine	3.646	79	270880	30.244	ng/u1	94
6) Benzaldehyde	6.911	77	174609	36.346	ng/u1	94
8) Phenol	6.981	94	309115	27.845	ng/u1	99
10) Bis(2-Chloroethyl)ether	7.199	93	243933	27.394	ng/u1	99
12) 2-Chlorophenol	7.328	128	227029	29.018	ng/u1	97
13) 2-Methylphenol	8.228	108	229813	27.599	ng/u1	99
14) 2,2'-oxybis(1-Chloropr...	8.305	45	303606	28.653	ng/u1	97
16) Acetophenone	8.599	105	390950	28.075	ng/u1	96
17) N-Nitroso-di-n-propyla...	8.581	70	211249	29.115	ng/u1	97
18) 4-Methylphenol	8.558	108	246052	27.757	ng/u1	98
19) Hexachloroethane	8.834	117	101293	28.663	ng/u1	93
22) Nitrobenzene	8.975	77	322496	29.578	ng/u1	98
23) Isophorone	9.505	82	595264	28.162	ng/u1	99
25) 2-Nitrophenol	9.687	139	135093	28.400	ng/u1	94
26) 2,4-Dimethylphenol	9.757	107	279928	27.292	ng/u1	97
27) Bis(2-Chloroethoxy)met...	9.987	93	338843	27.872	ng/u1	100
29) 2,4-Dichlorophenol	10.222	162	254825	28.902	ng/u1	99
30) Naphthalene	10.610	128	732637	28.065	ng/u1	100
32) 4-Chloroaniline	10.740	127	245227	21.689	ng/u1	100
33) Hexachlorobutadiene	10.887	225	219322	28.831	ng/u1	99
34) Caprolactam	11.540	113	77443	28.008	ng/u1	97
35) 4-Chloro-3-methylphenol	11.881	107	283301	28.942	ng/u1	97
36) 2-Methylnaphthalene	12.228	142	522111	27.489	ng/u1	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.451	142	526491	28.331	ng/u1	99
39) 1,2,4,5-Tetrachloroben...	12.598	216	399044	28.144	ng/u1	98
40) Hexachlorocyclopentadiene	12.569	237	232225	25.519	ng/u1	98
41) 2,4,6-Trichlorophenol	12.851	196	242512	27.794	ng/u1	99
42) 2,4,5-Trichlorophenol	12.934	196	263309	28.160	ng/u1	100
43) 1,1'-Biphenyl	13.251	154	748664	27.699	ng/u1	99
44) 2-Chloronaphthalene	13.293	162	604301	27.646	ng/u1	99
45) 2-Nitroaniline	13.516	65	191003	30.465	ng/u1	99
47) Dimethylphthalate	13.887	163	813260	28.309	ng/u1	100
48) 2,6-Dinitrotoluene	14.016	165	170158	29.953	ng/u1	98
50) Acenaphthylene	14.140	152	947965	27.932	ng/u1	99
51) 3-Nitroaniline	14.345	138	136145	27.476	ng/u1	99
52) Acenaphthene	14.487	153	631736	28.409	ng/u1	99
53) 2,4-Dinitrophenol	14.563	184	90984	24.426	ng/u1	96
55) 4-Nitrophenol	14.681	109	125718	26.227	ng/u1	96
56) Dibenzofuran	14.822	168	917961	28.833	ng/u1	99
57) 2,4-Dinitrotoluene	14.810	165	243312	30.526	ng/u1	96
58) 2,3,4,6-Tetrachlorophenol	15.057	232	245668	29.595	ng/u1	99
59) Diethylphthalate	15.251	149	785206	29.208	ng/u1	100
61) Fluorene	15.475	166	760282	29.155	ng/u1	99
62) 4-Chlorophenyl-phenyle...	15.469	204	455285	29.257	ng/u1	99
63) 4-Nitroaniline	15.516	138	129053	33.595	ng/u1	96
66) 4,6-Dinitro-2-methylph...	15.575	198	165802	27.464	ng/u1	97
67) N-Nitrosodiphenylamine	15.687	169	644836	28.353	ng/u1	99
68) 4-Bromophenyl-phenylether	16.369	248	296782	28.163	ng/u1	97
69) Hexachlorobenzene	16.481	284	339675	27.969	ng/u1	97
70) Atrazine	16.651	200	233176	23.520	ng/u1	98
71) Pentachlorophenol	16.834	266	205928	27.300	ng/u1	100
72) Phenanthrene	17.222	178	1220682	28.167	ng/u1	99
74) Anthracene	17.316	178	1225154	27.887	ng/u1	99
75) 1,2,3,4-Tetrachloroben...	13.210	216	397850	27.851	ng/uL	100
76) Pentachlorobenzene	14.740	250	394534	28.432	ng/uL	99
77) Carbazole	17.598	167	1025112	27.427	ng/u1	99
78) Di-n-butylphthalate	18.145	149	1261729	28.208	ng/u1	99
80) Fluoranthene	19.204	202	1510197	28.706	ng/u1	99
82) Pyrene	19.557	202	1529099	28.636	ng/u1	99
83) Butylbenzylphthalate	20.439	149	517677	28.258	ng/u1	94
84) 3,3'-Dichlorobenzidine	21.210	252	448641	28.948	ng/u1	99
85) Benzo(a)anthracene	21.274	228	1506350	28.349	ng/u1	99
86) Bis(2-ethylhexyl)phtha...	21.198	149	723602	27.517	ng/u1	100
87) Chrysene	21.327	228	1381653	28.446	ng/u1	99
89) Di-n-octyl phthalate	22.104	149	1162790	24.527	ng/u1	100
90) Benzo(b)fluoranthene	22.927	252	1381968	27.841	ng/u1	99
91) Benzo(k)fluoranthene	22.974	252	1345063	27.933	ng/u1	99
93) Benzo(a)pyrene	23.545	252	1265781	28.379	ng/u1	99
94) Indeno(1,2,3-cd)pyrene	26.092	276	1520720	28.829	ng/u1	99
95) Dibenzo(a,h)anthracene	26.110	278	1255471	28.750	ng/u1	100
96) Benzo(g,h,i)perylene	26.845	276	1203927	28.420	ng/u1	99

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

