

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072320\
 Data File : BP002841.D
 Acq On : 23 Jul 2020 15:37
 Operator : CG/JU
 Sample : L3380-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VNJ-241

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:13:11 PM

Quant Time: Jul 23 17:16:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	90058	20.00	ng	-0.01
21) Naphthalene-d8	10.31	136	389935	20.00	ng	-0.02
39) Acenaphthene-d10	14.20	164	186932	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	188837	20.00	ng	0.00
76) Chrysene-d12	21.08	240	320765	20.00	ng	0.00
86) Perylene-d12	23.26	264	542250	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	675576	126.57	ng	-0.01
7) Phenol-d6	6.75	99	933024	122.48	ng	0.00
23) Nitrobenzene-d5	8.69	82	608209	83.42	ng	-0.02
42) 2,4,6-Tribromophenol	15.72	330	206033	105.59	ng	0.00
45) 2-Fluorobiphenyl	12.81	172	1190507	98.52	ng	-0.01
79) Terphenyl-d14	19.57	244	811441	58.00	ng	0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.10	88	83030	35.679	ng	98
3) Pyridine	3.49	79	216343	31.195	ng	98
4) n-Nitrosodimethylamine	3.40	42	109916	42.071	ng	93
6) Aniline	6.88	93	239725	24.858	ng	99
8) 2-Chlorophenol	7.12	128	278873	46.958	ng	99
9) Benzaldehyde	6.69	77	219836	53.115	ng	97
10) Phenol	6.78	94	364615	46.724	ng	97
11) bis(2-Chloroethyl)ether	6.98	93	277380	42.735	ng	99
12) 1,3-Dichlorobenzene	7.43	146	297415	43.674	ng	97
13) 1,4-Dichlorobenzene	7.58	146	303493	44.408	ng	99
14) 1,2-Dichlorobenzene	7.89	146	291952	44.299	ng	99
15) Benzyl Alcohol	7.79	79	258445	50.698	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.08	45	397465	42.579	ng	99
17) 2-Methylphenol	8.01	107	245066	44.080	ng	99
18) Hexachloroethane	8.60	117	117538	48.557	ng	100
19) n-Nitroso-di-n-propylamine	8.35	70	206470	44.507	ng	98
20) 3+4-Methylphenols	8.34	107	320736	43.631	ng	# 89
22) Acetophenone	8.36	105	406880	42.327	ng	# 96
24) Nitrobenzene	8.73	77	334497	42.363	ng	99
25) Isophorone	9.26	82	619500	41.691	ng	99
26) 2-Nitrophenol	9.44	139	160611	47.095	ng	95
27) 2,4-Dimethylphenol	9.52	122	269203	48.621	ng	98
28) bis(2-Chloroethoxy)methane	9.74	93	382758	42.408	ng	99
29) 2,4-Dichlorophenol	9.99	162	281472	46.111	ng	98
30) 1,2,4-Trichlorobenzene	10.19	180	302160	43.216	ng	100
31) Naphthalene	10.36	128	885509	43.013	ng	100
32) Benzoic acid	9.72	122	183106	43.520	ng	98
33) 4-Chloroaniline	10.49	127	138990	15.756	ng	97
34) Hexachlorobutadiene	10.66	225	173507	42.934	ng	97
35) Caprolactam	11.26	113	103960	48.709	ng	# 91
36) 4-Chloro-3-methylphenol	11.65	107	296407	45.048	ng	99
37) 2-Methylnaphthalene	11.99	142	631979	43.503	ng	98
38) 1-Methylnaphthalene	12.21	142	587392	42.749	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.37	216	322883	56.203	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.35	237	203629	80.018	nq	98
43) 2,4,6-Trichlorophenol	12.63	196	203205	55.260	nq	97
44) 2,4,5-Trichlorophenol	12.72	196	225706	52.920	nq	# 95
46) 1,1'-Biphenyl	13.02	154	728622	52.283	nq	98
47) 2-Chloronaphthalene	13.06	162	563112	49.748	nq	97
48) 2-Nitroaniline	13.27	65	187938	53.714	nq	94
49) Acenaphthylene	13.92	152	836872	46.899	nq	99
50) Dimethylphthalate	13.67	163	833862	58.491	nq	98
51) 2,6-Dinitrotoluene	13.79	165	170304	55.800	nq	99
52) Acenaphthene	14.27	154	441097	41.362	nq	99
53) 3-Nitroaniline	14.12	138	103193	30.220	nq	# 97
54) 2,4-Dinitrophenol	14.35	184	127590	75.661	nq	96
55) Dibenzofuran	14.61	168	653227	38.192	nq	94
56) 4-Nitrophenol	14.47	139	202898	74.513	nq	# 71
57) 2,4-Dinitrotoluene	14.59	165	161940	40.289	nq	# 94
58) Fluorene	15.26	166	421633	33.316	nq	98
59) 2,3,4,6-Tetrachlorophenol	14.85	232	136122	40.489	nq	# 93
60) Diethylphthalate	15.05	149	480545	34.766	nq	98
61) 4-Chlorophenyl-phenylether	15.26	204	220285	32.759	nq	# 89
62) 4-Nitroaniline	15.29	138	82666m	23.401	nq	
63) Azobenzene	15.56	77	482426	33.945	nq	94
65) 4,6-Dinitro-2-methylphenol	15.37	198	74935	64.371	nq	# 55
66) n-Nitrosodiphenylamine	15.48	169	414573	74.319	nq	98
67) 4-Bromophenyl-phenylether	16.16	248	117199	56.976	nq	# 87
68) Hexachlorobenzene	16.30	284	126693	58.458	nq	# 84
69) Atrazine	16.46	200	106537	69.957	nq	84
70) Pentachlorophenol	16.65	266	119210	110.590	nq	99
71) Phenanthrene	17.02	178	431863	42.239	nq	# 90
72) Anthracene	17.12	178	448735	44.725	nq	# 95
73) Carbazole	17.39	167	378628	38.456	nq	# 87
74) Di-n-butylphthalate	17.96	149	470901	44.100	nq	# 96
75) Fluoranthene	19.02	202	547128	46.183	nq	94
77) Benzidine	19.19	184	235489	29.618	nq	# 84
78) Pyrene	19.36	202	592484	26.753	nq	98
80) Butylbenzylphthalate	20.24	149	306784	35.103	nq	90
81) Benzo(a)anthracene	21.07	228	899638	43.327	nq	99
82) 3,3'-Dichlorobenzidine	21.00	252	42982	6.250	nq	# 93
83) Chrysene	21.12	228	844418	42.723	nq	98
84) Bis(2-ethylhexyl)phthalate	21.00	149	554243	45.597	nq	100
85) Di-n-octyl phthalate	21.86	149	1235535	56.538	nq	98
87) Indeno(1,2,3-cd)pyrene	25.49	276	1833005	46.770	nq	98
88) Benzo(b)fluoranthene	22.62	252	1340911	39.824	nq	98
89) Benzo(k)fluoranthene	22.66	252	1335622	40.272	nq	100
90) Benzo(a)pyrene	23.17	252	1345566	42.171	nq	98
91) Dibenzo(a,h)anthracene	25.49	278	1517960	48.026	nq	99
92) Benzo(g,h,i)perylene	26.15	276	1618250	50.497	nq	99

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

