

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070920\
 Data File : BP002700.D
 Acq On : 09 Jul 2020 16:48
 Operator : CG/JU
 Sample : PB129999BSD
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB129999BSD

Manual Integrations
 APPROVED

mohammad
 7/10/2020 11:52:10 AM

Quant Time: Jul 09 17:29:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	68652	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	325208	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	235816	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	597092	20.00	ng	0.00
76) Chrysene-d12	21.09	240	777958	20.00	ng	0.00
86) Perylene-d12	23.28	264	886181	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.19	112	565689	139.51	ng	0.00
7) Phenol-d6	6.77	99	818427	144.61	ng	0.00
23) Nitrobenzene-d5	8.71	82	468921	74.29	ng	-0.01
42) 2,4,6-Tribromophenol	15.72	330	336688	117.39	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	1143692	71.38	ng	-0.01
79) Terphenyl-d14	19.57	244	2439008	67.67	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.13	88	54248	31.215	ng	100
3) Pyridine	3.52	79	170766	33.154	ng	98
4) n-Nitrosodimethylamine	3.43	42	79252	36.248	ng	# 86
6) Aniline	6.90	93	202784	28.522	ng	97
8) 2-Chlorophenol	7.14	128	214983	47.730	ng	99
9) Benzaldehyde	6.72	77	68468	20.777	ng	97
10) Phenol	6.79	94	303430	53.166	ng	95
11) bis(2-Chloroethyl)ether	7.00	93	204044	43.560	ng	99
12) 1,3-Dichlorobenzene	7.46	146	193582	37.245	ng	99
13) 1,4-Dichlorobenzene	7.60	146	199855	37.744	ng	98
14) 1,2-Dichlorobenzene	7.92	146	195890	38.433	ng	99
15) Benzyl Alcohol	7.81	79	188767	44.389	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.10	45	277570	42.530	ng	99
17) 2-Methylphenol	8.03	107	198444	46.454	ng	96
18) Hexachloroethane	8.63	117	69422	36.326	ng	98
19) n-Nitroso-di-n-propylamine	8.37	70	166238	44.769	ng	97
20) 3+4-Methylphenols	8.35	107	276002	47.923	ng	94
22) Acetophenone	8.38	105	313180	38.047	ng	# 97
24) Nitrobenzene	8.75	77	249263	37.311	ng	99
25) Isophorone	9.27	82	489431	38.690	ng	99
26) 2-Nitrophenol	9.46	139	113864	37.371	ng	94
27) 2,4-Dimethylphenol	9.55	122	210651	45.627	ng	97
28) bis(2-Chloroethoxy)methane	9.76	93	297708	40.969	ng	100
29) 2,4-Dichlorophenol	10.01	162	225046	42.295	ng	99
30) 1,2,4-Trichlorobenzene	10.21	180	207654	34.666	ng	100
31) Naphthalene	10.39	128	644883	37.459	ng	99
32) Benzoic acid	9.68	122	122472	36.353	ng	99
33) 4-Chloroaniline	10.50	127	174992	23.238	ng	99
34) Hexachlorobutadiene	10.69	225	110094	30.027	ng	97
35) Caprolactam	11.26	113	95229	53.242	ng	96
36) 4-Chloro-3-methylphenol	11.66	107	264401	45.338	ng	99
37) 2-Methylnaphthalene	12.01	142	498664	39.726	ng	97
38) 1-Methylnaphthalene	12.23	142	477353	40.378	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.39	216	251876	34.207	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.37	237	159251	43.543	nq	95
43) 2,4,6-Trichlorophenol	12.65	196	189005	37.935	nq	97
44) 2,4,5-Trichlorophenol	12.73	196	208343	36.326	nq	96
46) 1,1'-Biphenyl	13.04	154	644324	36.121	nq	98
47) 2-Chloronaphthalene	13.08	162	522862	36.009	nq	95
48) 2-Nitroaniline	13.29	65	190951	41.222	nq	96
49) Acenaphthylene	13.93	152	883233	38.650	nq	99
50) Dimethylphthalate	13.68	163	731780	39.510	nq	99
51) 2,6-Dinitrotoluene	13.80	165	164375	41.092	nq	95
52) Acenaphthene	14.28	154	520690	38.619	nq	100
53) 3-Nitroaniline	14.13	138	127374	29.244	nq	100
54) 2,4-Dinitrophenol	14.35	184	150624	61.877	nq	90
55) Dibenzofuran	14.62	168	858537	39.293	nq	98
56) 4-Nitrophenol	14.48	139	311608	90.305	nq	93
57) 2,4-Dinitrotoluene	14.60	165	243498	45.251	nq	97
58) Fluorene	15.27	166	698969	42.004	nq	99
59) 2,3,4,6-Tetrachlorophenol	14.86	232	190949	41.214	nq	97
60) Diethylphthalate	15.06	149	758057	41.160	nq	98
61) 4-Chlorophenyl-phenylether	15.27	204	346563	38.835	nq	92
62) 4-Nitroaniline	15.30	138	206191	47.365	nq	89
63) Azobenzene	15.57	77	768176	43.136	nq	97
65) 4,6-Dinitro-2-methylphenol	15.37	198	133224	35.419	nq	94
66) n-Nitrosodiphenylamine	15.49	169	659424	37.832	nq	98
67) 4-Bromophenyl-phenylether	16.17	248	223468	33.305	nq	96
68) Hexachlorobenzene	16.29	284	252517	35.129	nq	99
69) Atrazine	16.46	200	235141	42.042	nq	97
70) Pentachlorophenol	16.65	266	287031	70.207	nq	99
71) Phenanthrene	17.02	178	1284665	39.988	nq	100
72) Anthracene	17.11	178	1312475	41.055	nq	99
73) Carbazole	17.39	167	1325823	42.913	nq	99
74) Di-n-butylphthalate	17.96	149	1500681	41.243	nq	99
75) Fluoranthene	19.00	202	1721420	44.268	nq	98
77) Benzidine	19.19	184	816407	33.899	nq	99
78) Pyrene	19.36	202	1762345	33.281	nq	99
80) Butylbenzylphthalate	20.25	149	808421	34.970	nq	99
81) Benzo(a)anthracene	21.07	228	1920065	37.752	nq	98
82) 3,3'-Dichlorobenzidine	21.01	252	541694	29.091	nq	94
83) Chrysene	21.12	228	1829879	37.520	nq	100
84) Bis(2-ethylhexyl)phthalate	21.02	149	1195424	36.011	nq	98
85) Di-n-octyl phthalate	21.88	149	2266331	38.653	nq	99
87) Indeno(1,2,3-cd)pyrene	25.50	276	2809498	44.031	nq	96
88) Benzo(b)fluoranthene	22.63	252	2244842	39.699	nq	99
89) Benzo(k)fluoranthene	22.67	252	2005414m	38.016	nq	
90) Benzo(a)pyrene	23.19	252	2066712	39.828	nq	99
91) Dibenzo(a,h)anthracene	25.51	278	2317899	44.456	nq	99
92) Benzo(g,h,i)perylene	26.17	276	2370423	45.459	nq	97

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

