

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072420\
 Data File : BF120997.D
 Acq On : 24 Jul 2020 12:03
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 7/27/2020 12:19:25 PM

Quant Time: Jul 24 13:23:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	186666	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	690959	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	360265	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	683448	20.00	ng	0.00
76) Chrysene-d12	13.97	240	525889	20.00	ng	0.00
86) Perylene-d12	15.42	264	547958	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	819326	71.96	ng	0.00
7) Phenol-d6	6.45	99	1060653	72.11	ng	0.00
23) Nitrobenzene-d5	7.38	82	916823	76.78	ng	0.00
42) 2,4,6-Tribromophenol	10.64	330	312698	79.30	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1607778	66.63	ng	0.00
79) Terphenyl-d14	12.92	244	1815332	75.04	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.53	88	192919	36.813	ng	99
3) Pyridine	3.27	79	538720	38.644	ng	99
4) n-Nitrosodimethylamine	3.23	42	213091	35.747	ng	97
6) Aniline	6.48	93	698423	36.377	ng	98
8) 2-Chlorophenol	6.60	128	467575	37.275	ng	98
9) Benzaldehyde	6.36	77	289819	40.917	ng	99
10) Phenol	6.46	94	585346	36.178	ng	93
11) bis(2-Chloroethyl)ether	6.55	93	468379	37.773	ng	99
12) 1,3-Dichlorobenzene	6.76	146	503372	37.008	ng	98
13) 1,4-Dichlorobenzene	6.83	146	504993	37.337	ng	99
14) 1,2-Dichlorobenzene	6.99	146	467950	36.572	ng	99
15) Benzyl Alcohol	6.96	79	368781	35.454	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.09	45	680726	38.087	ng	99
17) 2-Methylphenol	7.07	107	415115	37.338	ng	99
18) Hexachloroethane	7.33	117	188774	38.562	ng	95
19) n-Nitroso-di-n-propylamine	7.23	70	301822	36.087	ng	98
20) 3+4-Methylphenols	7.22	107	454393	32.128	ng	95
22) Acetophenone	7.23	105	562463	36.271	ng	95
24) Nitrobenzene	7.40	77	493570	38.349	ng	99
25) Isophorone	7.64	82	904742	37.963	ng	99
26) 2-Nitrophenol	7.72	139	255586	41.813	ng	98
27) 2,4-Dimethylphenol	7.75	122	377090	37.996	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	569138	39.122	ng	100
29) 2,4-Dichlorophenol	7.96	162	397582	39.204	ng	99
30) 1,2,4-Trichlorobenzene	8.04	180	439662	39.076	ng	99
31) Naphthalene	8.12	128	1328244	37.476	ng	99
32) Benzoic acid	7.88	122	332514	44.633	ng	99
33) 4-Chloroaniline	8.17	127	593946	37.566	ng	99
34) Hexachlorobutadiene	8.24	225	256744	38.708	ng	100
35) Caprolactam	8.55	113	124889m	38.402	ng	
36) 4-Chloro-3-methylphenol	8.65	107	406480	39.082	ng	99
37) 2-Methylnaphthalene	8.81	142	885414	38.474	ng	99
38) 1-Methylnaphthalene	8.91	142	839396	38.512	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.97	216	413083	39.354	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.96	237	223718	38.564	ng	98
43) 2,4,6-Trichlorophenol	9.09	196	287660	38.923	ng	99
44) 2,4,5-Trichlorophenol	9.13	196	329088	39.255	ng	97
46) 1,1'-Biphenyl	9.27	154	1043940	37.988	ng	100
47) 2-Chloronaphthalene	9.30	162	856113	38.210	ng	99
48) 2-Nitroaniline	9.40	65	269985	39.874	ng	92
49) Acenaphthylene	9.72	152	1314692	37.771	ng	100
50) Dimethylphthalate	9.58	163	1006983	38.744	ng	99
51) 2,6-Dinitrotoluene	9.64	165	226808	40.618	ng	93
52) Acenaphthene	9.89	154	856405	38.700	ng	99
53) 3-Nitroaniline	9.81	138	273456	39.047	ng	96
54) 2,4-Dinitrophenol	9.91	184	116719	40.046	ng	87
55) Dibenzofuran	10.06	168	1207470	37.732	ng	100
56) 4-Nitrophenol	9.96	139	199315	37.466	ng	98
57) 2,4-Dinitrotoluene	10.04	165	304291	41.497	ng	95
58) Fluorene	10.40	166	846795	35.479	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.17	232	254579	39.884	ng	97
60) Diethylphthalate	10.27	149	1026274	39.038	ng	99
61) 4-Chlorophenyl-phenylether	10.39	204	440552	34.607	ng	99
62) 4-Nitroaniline	10.42	138	266623	39.084	ng	97
63) Azobenzene	10.55	77	947209	37.846	ng	97
65) 4,6-Dinitro-2-methylphenol	10.45	198	154226	39.711	ng	89
66) n-Nitrosodiphenylamine	10.51	169	831065	38.654	ng	99
67) 4-Bromophenyl-phenylether	10.88	248	305729	40.122	ng	97
68) Hexachlorobenzene	10.95	284	330125	40.050	ng	95
69) Atrazine	11.04	200	182803	34.332	ng	99
70) Pentachlorophenol	11.14	266	192934	40.857	ng	98
71) Phenanthrene	11.36	178	1295459	36.855	ng	100
72) Anthracene	11.41	178	1345210	38.113	ng	100
73) Carbazole	11.57	167	1316687	37.590	ng	100
74) Di-n-butylphthalate	11.90	149	1587831	39.282	ng	100
75) Fluoranthene	12.54	202	1455585	37.360	ng	99
77) Benzidine	12.67	184	511333	36.982	ng	99
78) Pyrene	12.77	202	1494200	40.188	ng	99
80) Butylbenzylphthalate	13.39	149	691064	41.694	ng	99
81) Benzo(a)anthracene	13.96	228	1194700	37.516	ng	98
82) 3,3'-Dichlorobenzidine	13.93	252	485874	40.441	ng	100
83) Chrysene	14.00	228	1292193	38.612	ng	99
84) Bis(2-ethylhexyl)phthalate	13.95	149	846155	41.400	ng	99
85) Di-n-octyl phthalate	14.56	149	1627325	40.020	ng	99
87) Indeno(1,2,3-cd)pyrene	16.84	276	1367902	35.895	ng	100
88) Benzo(b)fluoranthene	15.00	252	1373280	41.461	ng	100
89) Benzo(k)fluoranthene	15.03	252	1201721	39.783	ng	98
90) Benzo(a)pyrene	15.36	252	1217896	40.302	ng	100
91) Dibenzo(a,h)anthracene	16.87	278	1145041	36.784	ng	99
92) Benzo(g,h,i)perylene	17.27	276	1050332	34.221	ng	99

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

