

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060519\
 Data File : BM020748.D
 Acq On : 05 Jun 2019 22:13
 Operator : HP/JU
 Sample : IDOC-03-RC
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 IDOC-03-RC

Manual Integrations
 APPROVED

mohammad
 6/6/2019 11:22:06 AM

Quant Time: Jun 06 03:42:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 31 15:48:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	361058	20.00	ng	-0.01
21) Naphthalene-d8	10.61	136	1402176	20.00	ng	-0.02
39) Acenaphthene-d10	14.46	164	740730	20.00	ng	-0.02
64) Phenanthrene-d10	17.20	188	1404333	20.00	ng	-0.01
76) Chrysene-d12	21.38	240	1136335	20.00	ng	-0.02
87) Perylene-d12	23.67	264	1283434	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	2858293	139.26	ng	-0.01
7) Phenol-d6	6.99	99	3712778	122.85	ng	0.00
23) Nitrobenzene-d5	8.99	82	2227842	82.32	ng	-0.01
42) 2,4,6-Tribromophenol	15.95	330	1260732	137.37	ng	-0.01
45) 2-Fluorobiphenyl	13.08	172	4972360	89.24	ng	-0.01
79) Terphenyl-d14	19.83	244	5385414	79.27	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.35	88	328946	39.143	ng	95
3) Pyridine	3.74	79	944704	35.333	ng	97
4) n-Nitrosodimethylamine	3.66	42	427956	37.130	ng	92
6) Aniline	7.16	93	1328853	30.387	ng	98
8) 2-Chlorophenol	7.39	128	1109872	43.465	ng	98
9) Benzaldehyde	6.97	77	332031	15.187	ng	96
10) Phenol	7.02	94	1355761	41.626	ng	98
11) bis(2-Chloroethyl)ether	7.26	93	995321	38.345	ng	95
12) 1,3-Dichlorobenzene	7.71	146	1158028	39.460	ng	97
13) 1,4-Dichlorobenzene	7.86	146	1181713	39.610	ng	98
14) 1,2-Dichlorobenzene	8.17	146	1132212	38.997	ng	99
15) Benzyl Alcohol	8.06	79	978358	36.591	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.35	45	1340647	33.454	ng	99
17) 2-Methylphenol	8.26	107	902822	39.420	ng	98
18) Hexachloroethane	8.89	117	427710	37.445	ng	95
19) n-Nitroso-di-n-propylamine	8.63	70	812813	33.456	ng	97
20) 3+4-Methylphenols	8.59	107	1194848	38.345	ng	97
22) Acetophenone	8.65	105	1539322	43.215	ng	97
24) Nitrobenzene	9.03	77	1198515	43.381	ng	97
25) Isophorone	9.55	82	2113262	39.256	ng	99
26) 2-Nitrophenol	9.73	139	533231	51.787	ng	97
27) 2,4-Dimethylphenol	9.79	122	967381	50.138	ng	99
28) bis(2-Chloroethoxy)methane	10.03	93	1310681	40.268	ng	100
29) 2,4-Dichlorophenol	10.26	162	1007206	47.183	ng	98
30) 1,2,4-Trichlorobenzene	10.47	180	1110787	45.254	ng	98
31) Naphthalene	10.66	128	3163285	43.871	ng	99
32) Benzoic acid	9.93	122	541345m	44.942	ng	
33) 4-Chloroaniline	10.77	127	534308	15.922	ng	96
34) Hexachlorobutadiene	10.94	225	650958	44.200	ng	99
35) Caprolactam	11.57	113	272960m	36.807	ng	
36) 4-Chloro-3-methylphenol	11.89	107	1003435	40.430	ng	99
37) 2-Methylnaphthalene	12.27	142	2273239	42.534	ng	99
38) 1-Methylnaphthalene	12.49	142	2149452	42.237	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.64	216	1195654	49.256	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.62	237	1502099	103.465	ng	99
43) 2,4,6-Trichlorophenol	12.88	196	756300	48.403	ng	99
44) 2,4,5-Trichlorophenol	12.95	196	796525	49.378	ng	95
46) 1,1'-Biphenyl	13.29	154	2838104	45.877	ng	99
47) 2-Chloronaphthalene	13.33	162	2196582	43.994	ng	98
48) 2-Nitroaniline	13.54	65	600987	38.350	ng	92
49) Acenaphthylene	14.18	152	3420754	43.909	ng	100
50) Dimethylphthalate	13.92	163	2559534	40.445	ng	99
51) 2,6-Dinitrotoluene	14.04	165	557660	43.398	ng	89
52) Acenaphthene	14.52	154	1989418	42.851	ng	98
53) 3-Nitroaniline	14.37	138	351575	24.513	ng	95
54) 2,4-Dinitrophenol	14.57	184	466302	83.957	ng	91
55) Dibenzofuran	14.86	168	3111294	43.149	ng	97
56) 4-Nitrophenol	14.67	139	852712	80.679	ng	91
57) 2,4-Dinitrotoluene	14.83	165	718696	41.772	ng	94
58) Fluorene	15.51	166	2447357	41.273	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.08	232	660470	46.987	ng	95
60) Diethylphthalate	15.29	149	2449623	37.532	ng	99
61) 4-Chlorophenyl-phenylether	15.50	204	1307091	41.632	ng	94
62) 4-Nitroaniline	15.53	138	509191	33.475	ng	93
63) Azobenzene	15.80	77	2311242	36.937	ng	96
65) 4,6-Dinitro-2-methylphenol	15.59	198	286411	45.407	ng	93
66) n-Nitrosodiphenylamine	15.72	169	2066292	45.963	ng	99
67) 4-Bromophenyl-phenylether	16.40	248	756030	45.724	ng	95
68) Hexachlorobenzene	16.50	284	861764	44.633	ng	94
69) Atrazine	16.67	200	639604	47.670	ng	100
70) Pentachlorophenol	16.85	266	961646	104.403	ng	98
71) Phenanthrene	17.24	178	3422140	42.990	ng	99
72) Anthracene	17.33	178	3470139	43.555	ng	100
73) Carbazole	17.61	167	2950020	40.803	ng	99
74) Di-n-butylphthalate	18.17	149	3511472	37.414	ng	100
75) Fluoranthene	19.26	202	3556099	40.171	ng	99
77) Benzidine	19.45	184	1494350	43.400	ng	100
78) Pyrene	19.62	202	3568087	40.918	ng	100
80) Butylbenzylphthalate	20.52	149	1352411	36.699	ng	95
81) Benzo(a)anthracene	21.36	228	3276991	41.411	ng	99
82) 3,3'-Dichlorobenzidine	21.30	252	1060584	36.600	ng	99
83) Chrysene	21.41	228	3226344	42.892	ng	100
84) Bis(2-ethylhexyl)phthalate	21.29	149	1851436	34.568	ng	99
85) Di-n-octyl phthalate	22.19	149	3040474	35.757	ng	97
86) Indeno(1,2,3-cd)pyrene	25.99	276	4701606	61.096	ng	98
88) Benzo(b)fluoranthene	22.98	252	3697449	43.588	ng	99
89) Benzo(k)fluoranthene	23.03	252	3466222	41.813	ng	99
90) Benzo(a)pyrene	23.57	252	3598104	46.764	ng	99
91) Dibenzo(a,h)anthracene	26.00	278	3945514	52.855	ng	99
92) Benzo(g,h,i)perylene	26.70	276	3849001	55.512	ng	98

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

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