

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112618\
 Data File : BF110957.D
 Acq On : 26 Nov 2018 11:53
 Operator : JU/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

mohammad
 11/27/2018 11:36:51 AM

Quant Time: Nov 26 12:42:36 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 26 11:30:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	66957	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	288730	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	139416	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	265125	20.00	ng	0.00
76) Chrysene-d12	14.13	240	189315	20.00	ng	0.00
87) Perylene-d12	15.67	264	139683	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	493659	119.76	ng	0.00
7) Phenol-d6	6.57	99	618248	117.29	ng	0.00
23) Nitrobenzene-d5	7.52	82	597308	119.14	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	162881	115.95	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	1065799	109.18	ng	0.00
79) Terphenyl-d14	13.06	244	1122932	111.08	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.75	88	117991	57.448	ng	89
3) Pyridine	3.54	79	261204	58.918	ng	# 84
4) n-Nitrosodimethylamine	3.52	42	121419	63.830	ng	88
6) Aniline	6.61	93	390938	56.870	ng	# 56
8) 2-Chlorophenol	6.73	128	284353	58.843	ng	99
9) Benzaldehyde	6.50	77	147531	49.324	ng	99
10) Phenol	6.59	94	361327	59.115	ng	# 68
11) bis(2-Chloroethyl)ether	6.67	93	266364	58.149	ng	94
12) 1,3-Dichlorobenzene	6.88	146	311465	58.908	ng	98
13) 1,4-Dichlorobenzene	6.96	146	313290	58.350	ng	97
14) 1,2-Dichlorobenzene	7.11	146	296062	59.275	ng	100
15) Benzyl Alcohol	7.09	79	241465	58.536	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.21	45	414634	57.640	ng	74
17) 2-Methylphenol	7.19	107	224181	58.291	ng	# 86
18) Hexachloroethane	7.44	117	117530	59.294	ng	92
19) n-Nitroso-di-n-propylamine	7.36	70	203431	60.459	ng	96
20) 3+4-Methylphenols	7.35	107	291811	59.108	ng	95
22) Acetophenone	7.36	105	392028	58.259	ng	96
24) Nitrobenzene	7.53	77	305468	59.467	ng	93
25) Isophorone	7.76	82	493345	60.037	ng	98
26) 2-Nitrophenol	7.85	139	158196	61.733	ng	96
27) 2,4-Dimethylphenol	7.87	122	231240	59.251	ng	98
28) bis(2-Chloroethoxy)methane	7.96	93	329384	59.202	ng	99
29) 2,4-Dichlorophenol	8.09	162	241033	60.022	ng	99
30) 1,2,4-Trichlorobenzene	8.16	180	264894	58.506	ng	95
31) Naphthalene	8.25	128	780225	57.563	ng	100
32) Benzoic acid	8.05	122	176826	64.037	ng	97
33) 4-Chloroaniline	8.30	127	335723	54.682	ng	100
34) Hexachlorobutadiene	8.35	225	150496	58.194	ng	99
35) Caprolactam	8.70	113	82930	61.812	ng	90
36) 4-Chloro-3-methylphenol	8.78	107	265051	61.199	ng	97
37) 2-Methylnaphthalene	8.93	142	516219	58.097	ng	99
38) 1-Methylnaphthalene	9.03	142	500154	58.530	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.10	216	257108	58.261	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.07	237	37473	26.465	nq	98
43) 2,4,6-Trichlorophenol	9.22	196	160493	57.874	nq	96
44) 2,4,5-Trichlorophenol	9.27	196	164773	55.702	nq	94
46) 1,1'-Biphenyl	9.40	154	744349	57.233	nq	99
47) 2-Chloronaphthalene	9.43	162	544490	58.112	nq	99
48) 2-Nitroaniline	9.54	65	175041	60.624	nq	94
49) Acenaphthylene	9.85	152	767088	56.848	nq	99
50) Dimethylphthalate	9.70	163	591765	56.885	nq	99
51) 2,6-Dinitrotoluene	9.78	165	138421	60.488	nq	86
52) Acenaphthene	10.02	154	491306	57.614	nq	98
53) 3-Nitroaniline	9.96	138	157759	59.504	nq	86
54) 2,4-Dinitrophenol	10.07	184	44628	55.198	nq	89
55) Dibenzofuran	10.19	168	696180	55.801	nq	97
56) 4-Nitrophenol	10.12	139	86615	49.244	nq	98
57) 2,4-Dinitrotoluene	10.19	165	173271	58.234	nq	94
58) Fluorene	10.53	166	558439	58.274	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.31	232	129459	55.631	nq	93
60) Diethylphthalate	10.40	149	606942	58.974	nq	100
61) 4-Chlorophenyl-phenylether	10.52	204	289983	58.443	nq	96
62) 4-Nitroaniline	10.58	138	146605	54.913	nq	97
63) Azobenzene	10.68	77	593754	59.133	nq	98
65) 4,6-Dinitro-2-methylphenol	10.60	198	81510	61.002	nq	82
66) n-Nitrosodiphenylamine	10.65	169	514935	57.622	nq	97
67) 4-Bromophenyl-phenylether	11.01	248	169471	57.837	nq	# 89
68) Hexachlorobenzene	11.09	284	180386	58.391	nq	# 89
69) Atrazine	11.16	200	108324	39.644	nq	98
70) Pentachlorophenol	11.29	266	74936	45.459	nq	97
71) Phenanthrene	11.50	178	810966	57.094	nq	100
72) Anthracene	11.56	178	816374	56.976	nq	99
73) Carbazole	11.72	167	797714	57.971	nq	100
74) Di-n-butylphthalate	12.02	149	935357	57.965	nq	99
75) Fluoranthene	12.70	202	831617	58.398	nq	99
77) Benzidine	12.82	184	195261	37.505	nq	98
78) Pyrene	12.93	202	833542	57.183	nq	98
80) Butylbenzylphthalate	13.52	149	386114	61.542	nq	97
81) Benzo(a)anthracene	14.13	228	670892	59.289	nq	99
82) 3,3'-Dichlorobenzidine	14.08	252	207190	54.658	nq	98
83) Chrysene	14.16	228	644980	59.829	nq	98
84) Bis(2-ethylhexyl)phthalate	14.07	149	509686	65.486	nq	98
85) Di-n-octyl phthalate	14.69	149	782273	68.349	nq	98
86) Indeno(1,2,3-cd)pyrene	17.26	276	475498	61.466	nq	98
88) Benzo(b)fluoranthene	15.21	252	484795	58.015	nq	99
89) Benzo(k)fluoranthene	15.24	252	506708m	61.367	nq	
90) Benzo(a)pyrene	15.60	252	457893	60.310	nq	97
91) Dibenzo(a,h)anthracene	17.26	278	398612	62.041	nq	99
92) Benzo(g,h,i)perylene	17.74	276	392601	61.587	nq	98

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

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