

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112018\
 Data File : BF110887.D
 Acq On : 20 Nov 2018 11:32
 Operator : JU/SJ
 Sample : PB114904BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB114904BS

Manual Integrations
 APPROVED

Sohil
 11/21/2018 3:17:45 PM

Quant Time: Nov 20 11:59:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	63718	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	262488	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	123167	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	227416	20.00	ng	0.00
76) Chrysene-d12	14.13	240	166169	20.00	ng	0.00
87) Perylene-d12	15.66	264	120503	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	448714	114.39	ng	0.00
7) Phenol-d6	6.57	99	553674	110.38	ng	0.00
23) Nitrobenzene-d5	7.51	82	279032	61.22	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	137586	110.86	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	531518	61.63	ng	0.00
79) Terphenyl-d14	13.06	244	535992	60.41	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.86	88	59424	30.403	ng	87
3) Pyridine	3.64	79	158297	37.521	ng	# 83
4) n-Nitrosodimethylamine	3.59	42	79447	43.888	ng	84
6) Aniline	6.61	93	190347	29.098	ng	# 55
8) 2-Chlorophenol	6.73	128	173710	37.774	ng	99
9) Benzaldehyde	6.50	77	89810	31.522	ng	98
10) Phenol	6.59	94	209799	36.069	ng	# 64
11) bis(2-Chloroethyl)ether	6.67	93	156631	35.932	ng	94
12) 1,3-Dichlorobenzene	6.88	146	180928	35.959	ng	99
13) 1,4-Dichlorobenzene	6.96	146	186350	36.472	ng	99
14) 1,2-Dichlorobenzene	7.11	146	175134	36.846	ng	100
15) Benzyl Alcohol	7.09	79	143933	36.666	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.20	45	246950	36.074	ng	78
17) 2-Methylphenol	7.19	107	136052	37.174	ng	# 84
18) Hexachloroethane	7.45	117	70078	37.152	ng	96
19) n-Nitroso-di-n-propylamine	7.35	70	118089	36.879	ng	96
20) 3+4-Methylphenols	7.35	107	175531m	37.362	ng	
22) Acetophenone	7.35	105	238156	38.931	ng	# 94
24) Nitrobenzene	7.53	77	179361	38.408	ng	96
25) Isophorone	7.76	82	313542	41.971	ng	98
26) 2-Nitrophenol	7.85	139	88724	38.084	ng	94
27) 2,4-Dimethylphenol	7.87	122	154349	43.503	ng	99
28) bis(2-Chloroethoxy)methane	7.96	93	197054	38.959	ng	99
29) 2,4-Dichlorophenol	8.09	162	140776	38.561	ng	98
30) 1,2,4-Trichlorobenzene	8.16	180	154853	37.621	ng	97
31) Naphthalene	8.25	128	499748	40.556	ng	99
32) Benzoic acid	8.00	122	75167	29.943	ng	95
33) 4-Chloroaniline	8.30	127	131400	23.542	ng	99
34) Hexachlorobutadiene	8.35	225	86255	36.688	ng	97
35) Caprolactam	8.67	113	47088m	38.606	ng	
36) 4-Chloro-3-methylphenol	8.78	107	153228	38.917	ng	98
37) 2-Methylnaphthalene	8.94	142	340502	42.152	ng	99
38) 1-Methylnaphthalene	9.04	142	319562	41.135	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.10	216	145342	37.279	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.08	237	46870	35.524	nq	98
43) 2,4,6-Trichlorophenol	9.22	196	93919	38.335	nq	96
44) 2,4,5-Trichlorophenol	9.27	196	97189	37.189	nq	95
46) 1,1'-Biphenyl	9.40	154	411979	35.856	nq	99
47) 2-Chloronaphthalene	9.43	162	313035	37.817	nq	97
48) 2-Nitroaniline	9.53	65	97459	38.207	nq	96
49) Acenaphthylene	9.85	152	483554	40.564	nq	99
50) Dimethylphthalate	9.70	163	346985	37.755	nq	99
51) 2,6-Dinitrotoluene	9.77	165	78523	38.840	nq	87
52) Acenaphthene	10.02	154	289551	38.435	nq	98
53) 3-Nitroaniline	9.95	138	61778	26.376	nq	97
54) 2,4-Dinitrophenol	10.07	184	46732	59.318	nq	91
55) Dibenzofuran	10.19	168	417821	37.908	nq	98
56) 4-Nitrophenol	10.12	139	92606	59.596	nq	100
57) 2,4-Dinitrotoluene	10.19	165	100002	38.043	nq	98
58) Fluorene	10.54	166	343554	40.580	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.32	232	80958m	39.379	nq	
60) Diethylphthalate	10.40	149	351009	38.605	nq	99
61) 4-Chlorophenyl-phenylether	10.52	204	161702	36.889	nq	97
62) 4-Nitroaniline	10.57	138	77817	32.993	nq	96
63) Azobenzene	10.68	77	335401	37.810	nq	97
65) 4,6-Dinitro-2-methylphenol	10.60	198	36650	31.977	nq	78
66) n-Nitrosodiphenylamine	10.65	169	292757	38.192	nq	97
67) 4-Bromophenyl-phenylether	11.02	248	93356	37.143	nq	98
68) Hexachlorobenzene	11.09	284	98548	37.189	nq	# 89
69) Atrazine	11.17	200	94344	40.253	nq	98
70) Pentachlorophenol	11.29	266	77876	55.076	nq	97
71) Phenanthrene	11.50	178	492061	40.387	nq	100
72) Anthracene	11.56	178	515111	41.911	nq	98
73) Carbazole	11.72	167	431275	36.538	nq	99
74) Di-n-butylphthalate	12.02	149	533622	38.552	nq	98
75) Fluoranthene	12.70	202	495861	40.594	nq	100
77) Benzidine	12.82	184	201028	43.992	nq	96
78) Pyrene	12.93	202	505626	39.519	nq	98
80) Butylbenzylphthalate	13.53	149	217371	39.472	nq	97
81) Benzo(a)anthracene	14.12	228	402300	40.505	nq	99
82) 3,3'-Dichlorobenzidine	14.08	252	99196	29.814	nq	99
83) Chrysene	14.16	228	387036	40.903	nq	99
84) Bis(2-ethylhexyl)phthalate	14.07	149	284663	41.669	nq	96
85) Di-n-octyl phthalate	14.69	149	450893	44.883	nq	98
86) Indeno(1,2,3-cd)pyrene	17.25	276	294162	43.322	nq	98
88) Benzo(b)fluoranthene	15.21	252	295977	41.057	nq	100
89) Benzo(k)fluoranthene	15.24	252	301077	42.267	nq	99
90) Benzo(a)pyrene	15.60	252	279764	42.713	nq	97
91) Dibenzo(a,h)anthracene	17.25	278	248772	44.882	nq	99
92) Benzo(g,h,i)perylene	17.73	276	240050	43.650	nq	99

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

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