

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
 Data File : BF103867.D
 Acq On : 23 Mar 2018 9:30
 Operator : JU/SJ
 Sample : PB107555BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB107555BS

Manual Integrations
 APPROVED

Sohil
 3/26/2018 4:32:44 PM

Quant Time: Mar 23 14:01:30 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	144007	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	599770	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	281822	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	535720	20.00	ng	0.00
75) Chrysene-d12	14.17	240	463297	20.00	ng	0.00
86) Perylene-d12	15.72	264	449453	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	1051425	111.05	ng	0.01
7) Phenol-d6	6.60	99	1314116	113.45	ng	0.00
23) Nitrobenzene-d5	7.54	82	808241	93.98	ng	0.00
41) 2,4,6-Tribromophenol	10.82	330	365568	150.87	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1452383	82.21	ng	0.00
78) Terphenyl-d14	13.10	244	1547877	77.55	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.88	88	155255	33.302	ng	90
3) Pyridine	3.65	79	450158	35.105	ng	89
4) n-Nitrosodimethylamine	3.61	42	178496	37.753	ng	# 76
6) Aniline	6.64	93	364338	22.027	ng	96
8) 2-Chlorophenol	6.76	128	433032	43.660	ng	95
9) Benzaldehyde	6.53	77	119565	15.712	ng	94
10) Phenol	6.62	94	514205	41.776	ng	97
11) bis(2-Chloroethyl)ether	6.71	93	405007	39.859	ng	95
12) 1,3-Dichlorobenzene	6.92	146	443695	39.278	ng	99
13) 1,4-Dichlorobenzene	6.99	146	446398	39.326	ng	100
14) 1,2-Dichlorobenzene	7.15	146	418418	39.723	ng	99
15) Benzyl Alcohol	7.12	79	375853	41.878	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.25	45	537298	37.178	ng	93
17) 2-Methylphenol	7.22	107	363591	41.166	ng	99
18) Hexachloroethane	7.49	117	161403	40.423	ng	98
19) n-Nitroso-di-n-propylamine	7.39	70	283075	38.133	ng	92
20) 3+4-Methylphenols	7.37	107	446384	39.824	ng	# 84
22) Acetophenone	7.38	105	565937	36.715	ng	# 96
24) Nitrobenzene	7.56	77	440772	43.653	ng	98
25) Isophorone	7.79	82	826139	41.494	ng	99
26) 2-Nitrophenol	7.87	139	242050	60.054	ng	90
27) 2,4-Dimethylphenol	7.90	122	411645	48.262	ng	97
28) bis(2-Chloroethoxy)methane	8.00	93	524464	43.502	ng	100
29) 2,4-Dichlorophenol	8.11	162	365492	47.100	ng	99
30) 1,2,4-Trichlorobenzene	8.20	180	365693	40.631	ng	98
31) Naphthalene	8.28	128	1206472	39.821	ng	99
32) Benzoic acid	8.03	122	302689	49.821	ng	95
33) 4-Chloroaniline	8.32	127	99086	7.795	ng	96
34) Hexachlorobutadiene	8.39	225	191075	38.721	ng	98
35) Caprolactam	8.70	113	128739m	48.180	ng	
36) 4-Chloro-3-methylphenol	8.80	107	399241	46.669	ng	98
37) 2-Methylnaphthalene	8.97	142	813597	42.346	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.14	216	341788	42.511	ng	99
40) Hexachlorocyclopentadiene	9.12	237	383730	92.494	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.25	196	257489	50.069	nq	98
43) 2,4,5-Trichlorophenol	9.29	196	276911	47.707	nq	96
45) 1,1'-Biphenyl	9.43	154	959029	40.810	nq	99
46) 2-Chloronaphthalene	9.46	162	765866	41.032	nq	100
47) 2-Nitroaniline	9.56	65	242815	49.484	nq	99
48) Acenaphthylene	9.88	152	1185577	40.962	nq	99
49) Dimethylphthalate	9.73	163	939837	43.710	nq	100
50) 2,6-Dinitrotoluene	9.80	165	214032	51.171	nq	98
51) Acenaphthene	10.06	154	696015	40.499	nq	100
52) 3-Nitroaniline	9.97	138	124208	26.751	nq	98
53) 2,4-Dinitrophenol	10.08	184	210559	119.055	nq #	21
54) Dibenzofuran	10.23	168	1062910	43.304	nq	99
55) 4-Nitrophenol	10.13	139	388699	97.577	nq	95
56) 2,4-Dinitrotoluene	10.21	165	292593	54.552	nq #	90
57) Fluorene	10.57	166	831356	43.649	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.35	232	224539	54.594	nq	94
59) Diethylphthalate	10.43	149	914616	42.858	nq	98
60) 4-Chlorophenyl-phenylether	10.56	204	385864	44.330	nq	97
61) 4-Nitroaniline	10.59	138	238609	49.598	nq	93
62) Azobenzene	10.72	77	861237	39.694	nq	92
64) 4,6-Dinitro-2-methylphenol	10.62	198	142181	62.940	nq #	71
65) n-Nitrosodiphenylamine	10.68	169	751462	41.210	nq	99
66) 4-Bromophenyl-phenylether	11.05	248	236501	42.997	nq	90
67) Hexachlorobenzene	11.12	284	269124	42.869	nq #	89
68) Atrazine	11.20	200	251589	44.601	nq	99
69) Pentachlorophenol	11.32	266	295633	81.910	nq	98
70) Phenanthrene	11.54	178	1221162	41.671	nq	99
71) Anthracene	11.59	178	1263416	42.448	nq	100
72) Carbazole	11.75	167	1197317	41.388	nq	99
73) Di-n-butylphthalate	12.06	149	1445078	41.631	nq	100
74) Fluoranthene	12.73	202	1309857	43.386	nq	99
76) Benzidine	12.85	184	416640	21.085	nq	98
77) Pyrene	12.97	202	1322280	38.863	nq	100
79) Butylbenzylphthalate	13.57	149	669508	44.413	nq	99
80) Benzo(a)anthracene	14.16	228	1256999	42.862	nq	99
81) 3,3'-Dichlorobenzidine	14.12	252	239499	24.415	nq	99
82) Chrysene	14.20	228	1154264	41.289	nq	100
83) Bis(2-ethylhexyl)phthalate	14.13	149	909564	42.236	nq	98
84) Di-n-octyl phthalate	14.76	149	1606771	51.286	nq	96
85) Indeno(1,2,3-cd)pyrene	17.30	276	1071367	43.884	nq	98
87) Benzo(b)fluoranthene	15.26	252	1281941	45.146	nq	98
88) Benzo(k)fluoranthene	15.29	252	1058324	38.773	nq	98
89) Benzo(a)pyrene	15.66	252	1111750	43.167	nq	99
90) Dibenzo(a,h)anthracene	17.33	278	894210	40.097	nq	98
91) Benzo(g,h,i)perylene	17.79	276	871677	40.178	nq	98

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

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