

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113496.D
 Acq On : 2 Apr 2019 7:32
 Operator : JU/SJ
 Sample : PB118478BSD
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB118478BSD

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:50:02 PM

Quant Time: Apr 02 09:26:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	116007	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	450550	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	215990	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	411581	20.00	ng	0.00
76) Chrysene-d12	13.85	240	364735	20.00	ng	0.00
87) Perylene-d12	15.26	264	260962	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	345450	48.69	ng	0.01
7) Phenol-d6	6.35	99	442506	49.30	ng	0.00
23) Nitrobenzene-d5	7.27	82	381016	50.19	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	120364	45.11	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	765631	47.85	ng	0.00
79) Terphenyl-d14	12.79	244	765666	46.27	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.38	88	70743	19.653	ng	98
3) Pyridine	3.11	79	168235	18.958	ng	95
4) n-Nitrosodimethylamine	3.02	42	75813	19.218	ng	95
6) Aniline	6.37	93	209560	19.173	ng	95
8) 2-Chlorophenol	6.50	128	192201	23.329	ng	95
9) Benzaldehyde	6.25	77	125109	20.773	ng	97
10) Phenol	6.37	94	238894	23.315	ng	100
11) bis(2-Chloroethyl)ether	6.44	93	177137	22.553	ng	99
12) 1,3-Dichlorobenzene	6.65	146	205150	23.100	ng	98
13) 1,4-Dichlorobenzene	6.72	146	209727	24.459	ng	100
14) 1,2-Dichlorobenzene	6.87	146	197914	23.461	ng	99
15) Benzyl Alcohol	6.86	79	141115	23.833	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.99	45	282855	23.972	ng	90
17) 2-Methylphenol	6.97	107	150299	23.278	ng	98
18) Hexachloroethane	7.22	117	52902	17.177	ng	98
19) n-Nitroso-di-n-propylamine	7.12	70	127577	22.629	ng	98
20) 3+4-Methylphenols	7.13	107	195282	22.606	ng	92
22) Acetophenone	7.12	105	255680	24.884	ng	# 97
24) Nitrobenzene	7.29	77	191060	24.120	ng	96
25) Isophorone	7.53	82	349996	23.792	ng	98
26) 2-Nitrophenol	7.61	139	76476	18.265	ng	92
27) 2,4-Dimethylphenol	7.66	122	168994	28.059	ng	98
28) bis(2-Chloroethoxy)methane	7.74	93	227006	24.504	ng	98
29) 2,4-Dichlorophenol	7.86	162	162512	24.899	ng	99
30) 1,2,4-Trichlorobenzene	7.93	180	173726	24.671	ng	99
31) Naphthalene	8.01	128	551575	25.052	ng	99
32) Benzoic acid	7.75	122	50200m	10.354	ng	
33) 4-Chloroaniline	8.06	127	195274	22.744	ng	99
34) Hexachlorobutadiene	8.13	225	100819	24.089	ng	99
35) Caprolactam	8.42	113	47850	22.843	ng	100
36) 4-Chloro-3-methylphenol	8.56	107	161776	24.643	ng	99
37) 2-Methylnaphthalene	8.70	142	373677	27.200	ng	98
38) 1-Methylnaphthalene	8.80	142	348377	26.364	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.87	216	174533	25.173	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.86	237	1622	0.553	ng	94
43) 2,4,6-Trichlorophenol	8.99	196	104975	23.251	ng	97
44) 2,4,5-Trichlorophenol	9.03	196	114467	22.909	ng	98
46) 1,1'-Biphenyl	9.16	154	455629	25.257	ng	99
47) 2-Chloronaphthalene	9.19	162	345035	24.842	ng	98
48) 2-Nitroaniline	9.29	65	108779	24.956	ng	99
49) Acenaphthylene	9.60	152	553944	24.666	ng	99
50) Dimethylphthalate	9.46	163	398557	24.900	ng	100
51) 2,6-Dinitrotoluene	9.53	165	78808	21.824	ng	92
52) Acenaphthene	9.77	154	314119	24.837	ng	100
53) 3-Nitroaniline	9.70	138	105286	25.855	ng	97
54) 2,4-Dinitrophenol	9.83	184	1634	0.910	ng	# 65
55) Dibenzofuran	9.95	168	483561	25.433	ng	97
56) 4-Nitrophenol	9.88	139	96380	33.197	ng	89
57) 2,4-Dinitrotoluene	9.93	165	94736	20.630	ng	92
58) Fluorene	10.29	166	371186	25.131	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.07	232	92111	21.800	ng	97
60) Diethylphthalate	10.16	149	391284	24.919	ng	99
61) 4-Chlorophenyl-phenylether	10.27	204	188462	26.180	ng	94
62) 4-Nitroaniline	10.30	138	105980	25.277	ng	96
63) Azobenzene	10.43	77	348463	23.575	ng	95
65) 4,6-Dinitro-2-methylphenol	10.36	198	2573	1.143	ng	74
66) n-Nitrosodiphenylamine	10.39	169	333736	26.425	ng	100
67) 4-Bromophenyl-phenylether	10.76	248	113681	25.342	ng	98
68) Hexachlorobenzene	10.84	284	128740	24.725	ng	98
69) Atrazine	10.92	200	99427	24.942	ng	97
70) Pentachlorophenol	11.04	266	96996	35.754	ng	98
71) Phenanthrene	11.24	178	511324	25.017	ng	98
72) Anthracene	11.29	178	542818	26.146	ng	99
73) Carbazole	11.45	167	504715	25.477	ng	98
74) Di-n-butylphthalate	11.78	149	618532	26.924	ng	98
75) Fluoranthene	12.43	202	582593	25.214	ng	97
77) Benzidine	12.55	184	497976	35.653	ng	96
78) Pyrene	12.65	202	592655	23.448	ng	99
80) Butylbenzylphthalate	13.27	149	256683	23.044	ng	96
81) Benzo(a)anthracene	13.84	228	498315	23.868	ng	99
82) 3,3'-Dichlorobenzidine	13.80	252	213868	25.535	ng	98
83) Chrysene	13.87	228	513055	24.193	ng	100
84) Bis(2-ethylhexyl)phthalate	13.83	149	346484	25.374	ng	99
85) Di-n-octyl phthalate	14.43	149	627959	27.092	ng	100
86) Indeno(1,2,3-cd)pyrene	16.62	276	348274	19.136	ng	98
88) Benzo(b)fluoranthene	14.86	252	402742	24.844	ng	99
89) Benzo(k)fluoranthene	14.88	252	400030	28.100	ng	99
90) Benzo(a)pyrene	15.20	252	358760	24.963	ng	99
91) Dibenzo(a,h)anthracene	16.63	278	298102	23.989	ng	99
92) Benzo(g,h,i)perylene	17.03	276	286637	22.877	ng	98

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

