

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
 Data File : BF127167.D
 Acq On : 03 Mar 2022 13:10
 Operator : CG\JU
 Sample : PB143042BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB143042BS

Quant Time: Mar 03 16:59:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 01 23:57:09 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 03/04/2022
 Supervised By : Jagrut Upadhyay 03/04/2022

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 Upadhyay
 03/04/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	95763	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	385157	20.000	ng	# 0.00
39) Acenaphthene-d10	9.933	164	195593	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	365485	20.000	ng	# 0.00
76) Chrysene-d12	14.074	240	232572	20.000	ng	# 0.00
86) Perylene-d12	15.562	264	170499	20.000	ng	0.00

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System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	810341	130.786	ng	0.01
7) Phenol-d6	6.528	99	1022559	122.308	ng	0.00
23) Nitrobenzene-d5	7.463	82	714399	83.722	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	337605	149.914	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	1177694	88.641	ng	0.00
79) Terphenyl-d14	13.016	244	1471407	93.005	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.746	88	105610	37.249	ng	97
3) Pyridine	3.510	79	256464	34.487	ng	95
4) n-Nitrosodimethylamine	3.475	42	164979	45.288	ng	# 77
6) Aniline	6.557	93	366404	36.988	ng	96
8) 2-Chlorophenol	6.681	128	309166	46.505	ng	94
9) Benzaldehyde	6.445	77	181896	35.742	ng	99
10) Phenol	6.539	94	400471	47.406	ng	77
11) bis(2-Chloroethyl)ether	6.628	93	290873	43.900	ng	99
12) 1,3-Dichlorobenzene	6.834	146	320050	44.629	ng	98
13) 1,4-Dichlorobenzene	6.910	146	326421	44.898	ng	97
14) 1,2-Dichlorobenzene	7.063	146	308927	44.558	ng	97
15) Benzyl Alcohol	7.034	79	287580	40.830	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.163	45	384218	36.825	ng	88
17) 2-Methylphenol	7.145	107	259271	45.077	ng	# 90
18) Hexachloroethane	7.398	117	124788	44.779	ng	93
19) n-Nitroso-di-n-propyla...	7.304	70	224796	41.727	ng	98
20) 3+4-Methylphenols	7.298	107	327042	43.787	ng	92
22) Acetophenone	7.304	105	442876	43.672	ng	99
24) Nitrobenzene	7.481	77	354949	44.033	ng	97
25) Isophorone	7.716	82	617765	43.751	ng	# 96
26) 2-Nitrophenol	7.792	139	163906	47.757	ng	# 84
27) 2,4-Dimethylphenol	7.828	122	283027	50.198	ng	91
28) bis(2-Chloroethoxy)met...	7.922	93	349891	40.874	ng	99
29) 2,4-Dichlorophenol	8.034	162	238994	45.086	ng	97
30) 1,2,4-Trichlorobenzene	8.116	180	266058	44.895	ng	99
31) Naphthalene	8.198	128	887261	44.129	ng	99
32) Benzoic acid	7.963	122	177105	44.292	ng	91
33) 4-Chloroaniline	8.245	127	191643	24.216	ng	97
34) Hexachlorobutadiene	8.304	225	156967	45.368	ng	98
35) Caprolactam	8.628	113	81613m	44.089	ng	
36) 4-Chloro-3-methylphenol	8.733	107	306953	45.309	ng	95
37) 2-Methylnaphthalene	8.886	142	583307	44.711	ng	98
38) 1-Methylnaphthalene	8.986	142	548518	43.906	ng	97
40) 1,2,4,5-Tetrachloroben...	9.057	216	245605	47.702	ng	99
41) Hexachlorocyclopentadiene	9.039	237	281922	100.958	ng	96
43) 2,4,6-Trichlorophenol	9.169	196	169199	44.806	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.210	196	182397	45.900	ng	95
46) 1,1'-Biphenyl	9.357	154	694403	44.810	ng	96
47) 2-Chloronaphthalene	9.380	162	511417	44.504	ng	94
48) 2-Nitroaniline	9.480	65	199451	43.983	ng	97
49) Acenaphthylene	9.798	152	870735	46.647	ng	98
50) Dimethylphthalate	9.657	163	628498	44.954	ng	98
51) 2,6-Dinitrotoluene	9.722	165	139431	49.019	ng	# 84
52) Acenaphthene	9.969	154	609148m	50.178	ng	
53) 3-Nitroaniline	9.892	138	105060	30.216	ng	# 97
54) 2,4-Dinitrophenol	10.004	184	144364	95.888	ng	# 86
55) Dibenzofuran	10.145	168	738577	44.245	ng	93
56) 4-Nitrophenol	10.063	139	230137	89.616	ng	# 74
57) 2,4-Dinitrotoluene	10.127	165	188892	49.115	ng	# 92
58) Fluorene	10.486	166	604763	46.510	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.263	232	149848	47.565	ng	91
60) Diethylphthalate	10.357	149	653502	44.734	ng	99
61) 4-Chlorophenyl-phenyle...	10.475	204	278166	45.368	ng	98
62) 4-Nitroaniline	10.516	138	155548	45.303	ng	# 74
63) Azobenzene	10.639	77	667899	41.058	ng	93
65) 4,6-Dinitro-2-methylph...	10.539	198	95132	44.300	ng	95
66) n-Nitrosodiphenylamine	10.598	169	523027	43.642	ng	99
67) 4-Bromophenyl-phenylether	10.969	248	181566	44.463	ng	98
68) Hexachlorobenzene	11.033	284	203529	46.297	ng	92
69) Atrazine	11.127	200	179918	48.404	ng	96
70) Pentachlorophenol	11.233	266	209297	87.216	ng	97
71) Phenanthrene	11.451	178	914613	44.709	ng	99
72) Anthracene	11.504	178	932898	45.809	ng	99
73) Carbazole	11.663	167	800346	42.854	ng	99
74) Di-n-butylphthalate	11.980	149	1024442	42.669	ng	99
75) Fluoranthene	12.645	202	910330	46.867	ng	93
77) Benzidine	12.763	184	383120	60.617	ng	96
78) Pyrene	12.874	202	918879	45.477	ng	99
80) Butylbenzylphthalate	13.486	149	415191	43.405	ng	# 86
81) Benzo(a)anthracene	14.063	228	704429	44.928	ng	99
82) 3,3'-Dichlorobenzidine	14.021	252	154564	31.280	ng	# 97
83) Chrysene	14.104	228	657895	44.621	ng	99
84) Bis(2-ethylhexyl)phtha...	14.039	149	520610	41.455	ng	# 97
85) Di-n-octyl phthalate	14.657	149	759937	42.028	ng	95
87) Indeno(1,2,3-cd)pyrene	17.080	276	532627	47.642	ng	# 84
88) Benzo(b)fluoranthene	15.127	252	583204	50.929	ng	# 93
89) Benzo(k)fluoranthene	15.157	252	563992	51.063	ng	# 94
90) Benzo(a)pyrene	15.504	252	506148	56.587	ng	# 93
91) Dibenzo(a,h)anthracene	17.098	278	456493	49.370	ng	# 91
92) Benzo(g,h,i)perylene	17.545	276	457624	50.202	ng	# 85

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

