

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072822\
 Data File : BF129402.D
 Acq On : 28 Jul 2022 11:24
 Operator : CG\JU
 Sample : PB146497BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB146497BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 07/29/2022
 Supervised By : Jagrut Upadhyay 07/29/2022

Quant Time: Jul 29 02:06:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 17:24:17 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	241353	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	966283	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	499401	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	831656	20.000	ng	0.00
76) Chrysene-d12	14.068	240	559778	20.000	ng	0.00
86) Perylene-d12	15.557	264	434549	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	1831251	123.599	ng	0.03
7) Phenol-d6	6.534	99	2281914	122.533	ng	0.02
23) Nitrobenzene-d5	7.457	82	1524425	86.120	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	610961	127.247	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	2528560	78.325	ng	0.00
79) Terphenyl-d14	13.010	244	2655940	89.511	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.787	88	231729	34.665	ng	90
3) Pyridine	3.563	79	562087	30.278	ng	93
4) n-Nitrosodimethylamine	3.510	42	372565	45.703	ng	# 91
6) Aniline	6.551	93	813110	35.121	ng	95
8) 2-Chlorophenol	6.675	128	775541	47.912	ng	96
9) Benzaldehyde	6.440	77	208799	16.510	ng	97
10) Phenol	6.545	94	897268m	45.244	ng	
11) bis(2-Chloroethyl)ether	6.622	93	726564	46.196	ng	94
12) 1,3-Dichlorobenzene	6.828	146	780205	44.942	ng	97
13) 1,4-Dichlorobenzene	6.904	146	786265	44.533	ng	99
14) 1,2-Dichlorobenzene	7.057	146	756532	46.618	ng	100
15) Benzyl Alcohol	7.034	79	649780	48.109	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.157	45	982035	41.539	ng	84
17) 2-Methylphenol	7.145	107	601009	44.756	ng	96
18) Hexachloroethane	7.392	117	302488	45.500	ng	# 88
19) n-Nitroso-di-n-propyla...	7.304	70	498130	44.184	ng	93
20) 3+4-Methylphenols	7.298	107	705115	41.298	ng	# 83
22) Acetophenone	7.298	105	982886	43.690	ng	# 92
24) Nitrobenzene	7.475	77	785802	43.110	ng	95
25) Isophorone	7.710	82	1431828	45.126	ng	99
26) 2-Nitrophenol	7.787	139	409085	45.492	ng	96
27) 2,4-Dimethylphenol	7.828	122	696140m	51.609	ng	
28) bis(2-Chloroethoxy)met...	7.916	93	864453	46.965	ng	99
29) 2,4-Dichlorophenol	8.028	162	605960	43.525	ng	98
30) 1,2,4-Trichlorobenzene	8.110	180	630330	43.741	ng	97
31) Naphthalene	8.192	128	2062095	42.004	ng	99
32) Benzoic acid	7.957	122	351457	36.042	ng	97
33) 4-Chloroaniline	8.239	127	571321	28.346	ng	99
34) Hexachlorobutadiene	8.304	225	365553	44.626	ng	100
35) Caprolactam	8.634	113	197393m	43.610	ng	
36) 4-Chloro-3-methylphenol	8.734	107	664885	45.028	ng	94
37) 2-Methylnaphthalene	8.881	142	1366386	42.829	ng	98
38) 1-Methylnaphthalene	8.981	142	1281289	41.603	ng	98
40) 1,2,4,5-Tetrachloroben...	9.051	216	585713	44.666	ng	98
41) Hexachlorocyclopentadiene	9.034	237	624647	106.975	ng	99
43) 2,4,6-Trichlorophenol	9.163	196	411992	43.257	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.216	196	425794	41.110	ng	# 92
46) 1,1'-Biphenyl	9.351	154	1590655	43.644	ng	100
47) 2-Chloronaphthalene	9.375	162	1256457	42.645	ng	99
48) 2-Nitroaniline	9.481	65	423957	43.273	ng	# 83
49) Acenaphthylene	9.792	152	1980037	44.228	ng	99
50) Dimethylphthalate	9.657	163	1468588	42.598	ng	100
51) 2,6-Dinitrotoluene	9.722	165	340364	44.110	ng	91
52) Acenaphthene	9.963	154	1363970m	46.646	ng	
53) 3-Nitroaniline	9.892	138	285779	31.364	ng	# 90
54) 2,4-Dinitrophenol	9.998	184	347688	88.265	ng	94
55) Dibenzofuran	10.139	168	1708972	42.397	ng	97
56) 4-Nitrophenol	10.063	139	530467	78.914	ng	92
57) 2,4-Dinitrotoluene	10.128	165	444461	44.093	ng	95
58) Fluorene	10.480	166	1349066	41.829	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.257	232	328375	41.026	ng	92
60) Diethylphthalate	10.351	149	1410389	42.320	ng	99
61) 4-Chlorophenyl-phenyle...	10.469	204	610413	42.931	ng	92
62) 4-Nitroaniline	10.516	138	363873	40.256	ng	97
63) Azobenzene	10.633	77	1396867	41.094	ng	95
65) 4,6-Dinitro-2-methylph...	10.533	198	224826	42.830	ng	96
66) n-Nitrosodiphenylamine	10.592	169	1189183	42.937	ng	96
67) 4-Bromophenyl-phenylether	10.963	248	398610	43.770	ng	98
68) Hexachlorobenzene	11.027	284	446552	45.255	ng	99
69) Atrazine	11.122	200	365549	43.453	ng	97
70) Pentachlorophenol	11.227	266	427904	79.768	ng	97
71) Phenanthrene	11.445	178	1909799	42.068	ng	100
72) Anthracene	11.498	178	1936468	43.406	ng	99
73) Carbazole	11.657	167	1834235	43.676	ng	100
74) Di-n-butylphthalate	11.974	149	2112022	42.230	ng	100
75) Fluoranthene	12.639	202	1980874	43.064	ng	100
77) Benzidine	12.757	184	609247	30.808	ng	99
78) Pyrene	12.869	202	2023014	43.217	ng	100
80) Butylbenzylphthalate	13.474	149	894922	44.076	ng	99
81) Benzo(a)anthracene	14.057	228	1662417	43.475	ng	99
82) 3,3'-Dichlorobenzidine	14.016	252	461211	36.531	ng	# 98
83) Chrysene	14.098	228	1513443	41.996	ng	100
84) Bis(2-ethylhexyl)phtha...	14.027	149	1173091	43.886	ng	100
85) Di-n-octyl phthalate	14.645	149	1734217	45.084	ng	97
87) Indeno(1,2,3-cd)pyrene	17.068	276	1341886	45.856	ng	99
88) Benzo(b)fluoranthene	15.115	252	1410700	48.940	ng	98
89) Benzo(k)fluoranthene	15.151	252	1287058	45.564	ng	99
90) Benzo(a)pyrene	15.498	252	1269699	54.262	ng	100
91) Dibenzo(a,h)anthracene	17.092	278	1154332	48.466	ng	100
92) Benzo(g,h,i)perylene	17.539	276	1193964	49.059	ng	99

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

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