

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072622\
 Data File : BF129347.D
 Acq On : 26 Jul 2022 10:30
 Operator : CG\JU
 Sample : PB146423BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB146423BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 07/27/2022
 Supervised By : Jagrut Upadhyay 07/27/2022

Quant Time: Jul 27 00:38:36 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	6.881	152	280753	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	1138655	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	597013	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	1000730	20.000	ng	# 0.00
76) Chrysene-d12	14.068	240	644928	20.000	ng	0.00
86) Perylene-d12	15.551	264	526661	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	2075125	120.404	ng	0.02
7) Phenol-d6	6.522	99	2640727	121.900	ng	0.00
23) Nitrobenzene-d5	7.451	82	1742374	83.532	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	710912	123.856	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	2881248	74.657	ng	0.00
79) Terphenyl-d14	13.004	244	3053295	89.317	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.763	88	273526	35.175	ng	89
3) Pyridine	3.540	79	714380	33.082	ng	88
4) n-Nitrosodimethylamine	3.487	42	424064	44.720	ng	# 88
6) Aniline	6.551	93	1003280	37.254	ng	99
8) 2-Chlorophenol	6.669	128	902579	47.935	ng	98
9) Benzaldehyde	6.434	77	335029	22.773	ng	98
10) Phenol	6.534	94	1055826m	45.768	ng	
11) bis(2-Chloroethyl)ether	6.622	93	855000	46.734	ng	93
12) 1,3-Dichlorobenzene	6.828	146	899177	44.526	ng	97
13) 1,4-Dichlorobenzene	6.904	146	916383	44.619	ng	99
14) 1,2-Dichlorobenzene	7.057	146	877744	46.496	ng	99
15) Benzyl Alcohol	7.028	79	729348	46.422	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.157	45	1231411	44.777	ng	92
17) 2-Methylphenol	7.139	107	696916	44.615	ng	99
18) Hexachloroethane	7.398	117	349769	45.229	ng	93
19) n-Nitroso-di-n-propyla...	7.298	70	589744	44.969	ng	95
20) 3+4-Methylphenols	7.292	107	822133	41.394	ng	# 89
22) Acetophenone	7.298	105	1138057	42.929	ng	# 99
24) Nitrobenzene	7.469	77	914114	42.557	ng	97
25) Isophorone	7.710	82	1704459	45.586	ng	98
26) 2-Nitrophenol	7.787	139	479752	45.275	ng	92
27) 2,4-Dimethylphenol	7.822	122	816132	51.346	ng	97
28) bis(2-Chloroethoxy)met...	7.916	93	1048693	48.350	ng	100
29) 2,4-Dichlorophenol	8.028	162	717326	43.724	ng	98
30) 1,2,4-Trichlorobenzene	8.110	180	744677	43.853	ng	95
31) Naphthalene	8.192	128	2449940	42.349	ng	99
32) Benzoic acid	7.951	122	518260	45.102	ng	98
33) 4-Chloroaniline	8.239	127	690965	29.092	ng	97
34) Hexachlorobutadiene	8.304	225	418125	43.317	ng	100
35) Caprolactam	8.622	113	241900m	45.353	ng	
36) 4-Chloro-3-methylphenol	8.722	107	776782	44.642	ng	96
37) 2-Methylnaphthalene	8.881	142	1594551	42.414	ng	98
38) 1-Methylnaphthalene	8.981	142	1532843	42.237	ng	97
40) 1,2,4,5-Tetrachloroben...	9.051	216	671201	42.816	ng	# 98
41) Hexachlorocyclopentadiene	9.033	237	746358	106.920	ng	99
43) 2,4,6-Trichlorophenol	9.163	196	489797	43.018	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.210	196	527039	42.565	ng	# 93
46) 1,1'-Biphenyl	9.351	154	1845196	42.350	ng	100
47) 2-Chloronaphthalene	9.375	162	1479207	41.997	ng	99
48) 2-Nitroaniline	9.475	65	497363	42.466	ng	89
49) Acenaphthylene	9.792	152	2355160	44.006	ng	99
50) Dimethylphthalate	9.651	163	1744488	42.328	ng	99
51) 2,6-Dinitrotoluene	9.716	165	408331	44.266	ng	100
52) Acenaphthene	9.963	154	1602140m	45.833	ng	
53) 3-Nitroaniline	9.886	138	341528	31.354	ng	# 92
54) 2,4-Dinitrophenol	9.998	184	434787	92.330	ng	# 87
55) Dibenzofuran	10.139	168	2008018	41.671	ng	97
56) 4-Nitrophenol	10.051	139	682995	84.992	ng	96
57) 2,4-Dinitrotoluene	10.122	165	536784	44.545	ng	98
58) Fluorene	10.480	166	1555237	40.337	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.251	232	407618	42.600	ng	93
60) Diethylphthalate	10.351	149	1682346	42.227	ng	99
61) 4-Chlorophenyl-phenyle...	10.469	204	714212	42.018	ng	90
62) 4-Nitroaniline	10.510	138	439362	40.661	ng	93
63) Azobenzene	10.633	77	1681419	41.377	ng	95
65) 4,6-Dinitro-2-methylph...	10.533	198	279735	44.287	ng	87
66) n-Nitrosodiphenylamine	10.592	169	1415381	42.470	ng	96
67) 4-Bromophenyl-phenylether	10.963	248	483405	44.113	ng	96
68) Hexachlorobenzene	11.027	284	526328	44.327	ng	99
69) Atrazine	11.116	200	446166	44.076	ng	97
70) Pentachlorophenol	11.222	266	563716	87.332	ng	98
71) Phenanthrene	11.445	178	2266083	41.483	ng	99
72) Anthracene	11.498	178	2280227	42.476	ng	99
73) Carbazole	11.651	167	2178390	43.107	ng	100
74) Di-n-butylphthalate	11.974	149	2547811	42.336	ng	99
75) Fluoranthene	12.633	202	2319992	41.915	ng	97
77) Benzidine	12.757	184	772737	33.916	ng	99
78) Pyrene	12.869	202	2335377	43.303	ng	99
80) Butylbenzylphthalate	13.474	149	1062756	45.432	ng	98
81) Benzo(a)anthracene	14.057	228	1894822	43.011	ng	100
82) 3,3'-Dichlorobenzidine	14.016	252	516279	35.494	ng	98
83) Chrysene	14.092	228	1715656	41.321	ng	99
84) Bis(2-ethylhexyl)phtha...	14.027	149	1409851	45.779	ng	99
85) Di-n-octyl phthalate	14.645	149	2059334	46.468	ng	98
87) Indeno(1,2,3-cd)pyrene	17.068	276	1588751	44.797	ng	98
88) Benzo(b)fluoranthene	15.115	252	1572733	45.019	ng	98
89) Benzo(k)fluoranthene	15.145	252	1567851	45.796	ng	98
90) Benzo(a)pyrene	15.492	252	1473315	51.952	ng	97
91) Dibenzo(a,h)anthracene	17.086	278	1371500	47.513	ng	99
92) Benzo(g,h,i)perylene	17.527	276	1408591	47.755	ng	98

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

