

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
 Data File : BF124999.D
 Acq On : 9 Aug 2021 11:42
 Operator : JU/CG
 Sample : PB138219BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB138219BSD

Manual Integrations
 APPROVED

mohammad
 8/10/2021 4:34:26 PM

Quant Time: Aug 09 14:32:16 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 09 11:03:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	7652	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	30251	20.000	ng	# 0.00
39) Acenaphthene-d10	9.863	164	16558	20.000	ng	0.00
64) Phenanthrene-d10	11.345	188	28799	20.000	ng	0.00
76) Chrysene-d12	13.986	240	19403	20.000	ng	# 0.00
86) Perylene-d12	15.439	264	14937	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	44032	94.262	ng	0.01
7) Phenol-d6	6.457	99	53692	91.156	ng	0.00
23) Nitrobenzene-d5	7.392	82	54853	94.857	ng	0.00
42) 2,4,6-Tribromophenol	10.651	330	14414	97.439	ng	0.00
45) 2-Fluorobiphenyl	9.180	172	109021	96.342	ng	0.00
79) Terphenyl-d14	12.933	244	127312	110.680	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.675	88	6726	38.487	ng	# 100
3) Pyridine	3.416	79	15060	37.086	ng	94
4) n-Nitrosodimethylamine	3.381	42	14194	49.221	ng	# 84
6) Aniline	6.486	93	29444	47.324	ng	# 97
8) 2-Chlorophenol	6.610	128	23750	46.351	ng	92
9) Benzaldehyde	6.375	77	14357	44.802	ng	98
10) Phenol	6.475	94	28804	47.395	ng	82
11) bis(2-Chloroethyl)ether	6.557	93	22093	48.458	ng	92
12) 1,3-Dichlorobenzene	6.763	146	26182	46.946	ng	97
13) 1,4-Dichlorobenzene	6.839	146	26766	48.009	ng	98
14) 1,2-Dichlorobenzene	6.992	146	24728	47.328	ng	96
15) Benzyl Alcohol	6.969	79	20535	48.332	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.098	45	36472	49.931	ng	91
17) 2-Methylphenol	7.080	107	18817	49.060	ng	95
18) Hexachloroethane	7.333	117	10784	49.405	ng	# 86
19) n-Nitroso-di-n-propyla...	7.239	70	17345	49.629	ng	94
20) 3+4-Methylphenols	7.233	107	25093	49.283	ng	# 81
22) Acetophenone	7.233	105	35180	47.485	ng	# 86
24) Nitrobenzene	7.410	77	24897	44.494	ng	94
25) Isophorone	7.651	82	44043	44.624	ng	94
26) 2-Nitrophenol	7.722	139	13252	46.638	ng	# 82
27) 2,4-Dimethylphenol	7.763	122	21906	54.385	ng	89
28) bis(2-Chloroethoxy)met...	7.857	93	28699	51.091	ng	97
29) 2,4-Dichlorophenol	7.963	162	20177	44.961	ng	94
30) 1,2,4-Trichlorobenzene	8.045	180	22908	45.550	ng	97
31) Naphthalene	8.127	128	71189	45.803	ng	98
32) Benzoic acid	7.904	122	11974	49.823	ng	89
33) 4-Chloroaniline	8.175	127	23711	40.937	ng	# 93
34) Hexachlorobutadiene	8.239	225	14069	46.845	ng	95
35) Caprolactam	8.563	113	5599m	46.983	ng	
36) 4-Chloro-3-methylphenol	8.663	107	22248	47.394	ng	89
37) 2-Methylnaphthalene	8.816	142	49574	47.439	ng	98
38) 1-Methylnaphthalene	8.916	142	43515	44.503	ng	97
40) 1,2,4,5-Tetrachloroben...	8.986	216	23084	48.570	ng	97
41) Hexachlorocyclopentadiene	8.969	237	22733	122.618	ng	99
43) 2,4,6-Trichlorophenol	9.098	196	15467	48.887	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.139	196	15302	47.778	ng	97
46) 1,1'-Biphenyl	9.280	154	56446	45.746	ng	98
47) 2-Chloronaphthalene	9.310	162	46369	47.235	ng	99
48) 2-Nitroaniline	9.410	65	13808	47.884	ng	98
49) Acenaphthylene	9.727	152	71966	48.262	ng	99
50) Dimethylphthalate	9.586	163	54690	48.740	ng	# 95
51) 2,6-Dinitrotoluene	9.651	165	11788	47.179	ng	99
52) Acenaphthene	9.898	154	41604	46.023	ng	99
53) 3-Nitroaniline	9.822	138	9842	37.399	ng	# 85
54) 2,4-Dinitrophenol	9.933	184	11893	92.929	ng	85
55) Dibenzofuran	10.069	168	64552	45.692	ng	99
56) 4-Nitrophenol	9.992	139	17955	103.503	ng	91
57) 2,4-Dinitrotoluene	10.057	165	16214	47.797	ng	# 99
58) Fluorene	10.410	166	50860	46.950	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.186	232	12023	50.034	ng	# 92
60) Diethylphthalate	10.286	149	56518	49.083	ng	99
61) 4-Chlorophenyl-phenyle...	10.398	204	25340	49.534	ng	97
62) 4-Nitroaniline	10.439	138	12047	47.150	ng	91
63) Azobenzene	10.563	77	53204	47.249	ng	93
65) 4,6-Dinitro-2-methylph...	10.463	198	7815	43.316	ng	99
66) n-Nitrosodiphenylamine	10.521	169	44903	48.106	ng	97
67) 4-Bromophenyl-phenylether	10.892	248	15524	49.215	ng	93
68) Hexachlorobenzene	10.957	284	16748	48.804	ng	# 90
69) Atrazine	11.051	200	14942	53.273	ng	96
70) Pentachlorophenol	11.157	266	16228	99.144	ng	95
71) Phenanthrene	11.374	178	74865	47.922	ng	99
72) Anthracene	11.427	178	77531	49.471	ng	99
73) Carbazole	11.580	167	64384	44.798	ng	99
74) Di-n-butylphthalate	11.904	149	93497	51.188	ng	100
75) Fluoranthene	12.563	202	76812	47.528	ng	97
77) Benzidine	12.686	184	50127	96.553	ng	98
78) Pyrene	12.792	202	77238	50.538	ng	98
80) Butylbenzylphthalate	13.404	149	35914	52.804	ng	97
81) Benzo(a)anthracene	13.974	228	60698	48.004	ng	99
82) 3,3'-Dichlorobenzidine	13.939	252	17788	53.361	ng	# 99
83) Chrysene	14.015	228	57922	47.386	ng	100
84) Bis(2-ethylhexyl)phtha...	13.957	149	52540	55.687	ng	99
85) Di-n-octyl phthalate	14.568	149	82563	55.467	ng	98
87) Indeno(1,2,3-cd)pyrene	16.892	276	44078	49.554	ng	93
88) Benzo(b)fluoranthene	15.015	252	52430	49.990	ng	98
89) Benzo(k)fluoranthene	15.045	252	45019m	47.186	ng	
90) Benzo(a)pyrene	15.380	252	44374	49.685	ng	98
91) Dibenzo(a,h)anthracene	16.909	278	37049	47.372	ng	96
92) Benzo(g,h,i)perylene	17.333	276	36544	48.712	ng	# 89

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

