

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102521\
 Data File : BF125915.D
 Acq On : 25 Oct 2021 13:52
 Operator : JU/CG
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 10/26/2021 11:33:12 AM

Quant Time: Oct 25 17:25:17 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 25 17:23:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	140989	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	553650	20.000	ng	# 0.00
39) Acenaphthene-d10	9.904	164	307287	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	550830	20.000	ng	0.00
76) Chrysene-d12	14.027	240	368025	20.000	ng	# 0.00
86) Perylene-d12	15.492	264	263476	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	221776	25.697	ng	0.00
7) Phenol-d6	6.481	99	292020	26.284	ng	-0.01
23) Nitrobenzene-d5	7.428	82	219492	24.643	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	55374	18.252	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	440698	24.737	ng	0.00
79) Terphenyl-d14	12.974	244	472857	19.572	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.634	88	51855	14.111	ng	# 100
3) Pyridine	3.398	79	129912	13.665	ng	# 85
4) n-Nitrosodimethylamine	3.310	42	83581	24.424	ng	# 77
6) Aniline	6.528	93	168740	13.192	ng	# 92
8) 2-Chlorophenol	6.651	128	112487	11.900	ng	82
9) Benzaldehyde	6.416	77	97634	16.193	ng	93
10) Phenol	6.492	94	148846	13.777	ng	91
11) bis(2-Chloroethyl)ether	6.598	93	116426	13.300	ng	# 81
12) 1,3-Dichlorobenzene	6.810	146	123642	11.990	ng	97
13) 1,4-Dichlorobenzene	6.886	146	123058	11.701	ng	96
14) 1,2-Dichlorobenzene	7.039	146	118335	12.015	ng	97
15) Benzyl Alcohol	7.004	79	109630	14.889	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.145	45	274465	24.516	ng	94
17) 2-Methylphenol	7.110	107	91151	12.113	ng	92
18) Hexachloroethane	7.386	117	45833	12.656	ng	93
19) n-Nitroso-di-n-propyla...	7.275	70	82099	13.998	ng	# 93
20) 3+4-Methylphenols	7.263	107	122671	13.218	ng	88
22) Acetophenone	7.275	105	164264	13.497	ng	# 87
24) Nitrobenzene	7.445	77	118681	13.749	ng	88
25) Isophorone	7.686	82	218139	13.824	ng	# 87
26) 2-Nitrophenol	7.763	139	29647	6.792	ng	# 66
27) 2,4-Dimethylphenol	7.798	122	90227	12.413	ng	90
28) bis(2-Chloroethoxy)met...	7.898	93	145118	13.249	ng	# 96
29) 2,4-Dichlorophenol	7.998	162	84761	10.824	ng	93
30) 1,2,4-Trichlorobenzene	8.092	180	103085	12.005	ng	95
31) Naphthalene	8.175	128	323892	12.099	ng	98
32) Benzoic acid	7.863	122	38201	7.474	ng	94
33) 4-Chloroaniline	8.216	127	133256	11.895	ng	# 89
34) Hexachlorobutadiene	8.298	225	63882	13.318	ng	98
35) Caprolactam	8.557	113	26949	11.911	ng	# 42
36) 4-Chloro-3-methylphenol	8.686	107	97049	12.686	ng	89
37) 2-Methylnaphthalene	8.863	142	208546	12.029	ng	90
38) 1-Methylnaphthalene	8.963	142	202499	12.037	ng	90
40) 1,2,4,5-Tetrachloroben...	9.028	216	96359	11.684	ng	98
41) Hexachlorocyclopentadiene	9.022	237	46877	11.581	ng	98
43) 2,4,6-Trichlorophenol	9.133	196	61629	10.310	ng	92

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44) 2,4,5-Trichlorophenol	9.169	196	65018	10.230	ng	94
46) 1,1'-Biphenyl	9.328	154	258526	11.443	ng	97
47) 2-Chloronaphthalene	9.351	162	194960	10.628	ng	97
48) 2-Nitroaniline	9.439	65	54930	12.473	ng #	68
49) Acenaphthylene	9.763	152	297159	11.062	ng	98
50) Dimethylphthalate	9.622	163	213197	10.539	ng	97
51) 2,6-Dinitrotoluene	9.680	165	35957	8.794	ng #	67
52) Acenaphthene	9.939	154	187671	11.182	ng	99
53) 3-Nitroaniline	9.845	138	43143	8.924	ng #	63
54) 2,4-Dinitrophenol	9.945	184	8145	4.693	ng #	1
55) Dibenzofuran	10.110	168	288915	11.491	ng	92
56) 4-Nitrophenol	9.992	139	33433	9.277	ng #	74
57) 2,4-Dinitrotoluene	10.080	165	41847	8.094	ng #	94
58) Fluorene	10.451	166	213361	11.587	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.222	232	52197	10.880	ng #	89
60) Diethylphthalate	10.327	149	216952	11.240	ng	98
61) 4-Chlorophenyl-phenyle...	10.445	204	103786	11.701	ng	88
62) 4-Nitroaniline	10.451	138	38762	8.676	ng	91
63) Azobenzene	10.604	77	249606	13.758	ng	96
65) 4,6-Dinitro-2-methylph...	10.486	198	13737	5.000	ng	94
66) n-Nitrosodiphenylamine	10.557	169	187346	9.970	ng	99
67) 4-Bromophenyl-phenylether	10.933	248	63799	10.458	ng	95
68) Hexachlorobenzene	10.998	284	68595	9.922	ng	95
69) Atrazine	11.086	200	56813	10.417	ng	89
70) Pentachlorophenol	11.186	266	34730	9.398	ng	92
71) Phenanthrene	11.410	178	328671	11.152	ng	97
72) Anthracene	11.463	178	326806	11.102	ng	99
73) Carbazole	11.616	167	314598	11.484	ng	97
74) Di-n-butylphthalate	11.957	149	328686	10.868	ng	99
75) Fluoranthene	12.598	202	337455	12.690	ng	98
77) Benzidine	12.721	184	123838	9.195	ng	99
78) Pyrene	12.827	202	347995	9.307	ng	97
80) Butylbenzylphthalate	13.457	149	103610	8.645	ng #	83
81) Benzo(a)anthracene	14.015	228	270768	11.340	ng	100
82) 3,3'-Dichlorobenzidine	13.980	252	81523	10.806	ng #	96
83) Chrysene	14.051	228	269249	11.409	ng	98
84) Bis(2-ethylhexyl)phtha...	14.021	149	120974	9.277	ng #	96
85) Di-n-octyl phthalate	14.633	149	165642	7.624	ng #	90
87) Indeno(1,2,3-cd)pyrene	16.951	276	165918	8.251	ng	94
88) Benzo(b)fluoranthene	15.062	252	192738m	12.957	ng	
89) Benzo(k)fluoranthene	15.092	252	192598	12.611	ng	98
90) Benzo(a)pyrene	15.427	252	146751	11.140	ng	97
91) Dibenzo(a,h)anthracene	16.974	278	137972	8.318	ng	93
92) Benzo(g,h,i)perylene	17.392	276	139319	8.224	ng #	88

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

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Abundance

TIC: BF125915.D\data.ms

Time-->