

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042721\
 Data File : BF123746.D
 Acq On : 27 Apr 2021 14:32
 Operator : JU/CG
 Sample : SSTDIC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDIC050

Manual Integrations
 APPROVED

mohammad
 4/28/2021 10:59:58 AM

Quant Time: Apr 27 14:53:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 27 14:29:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.810	152	269387	20.00	ng	0.00
21) Naphthalene-d8	8.098	136	979303	20.00	ng	0.00
39) Acenaphthene-d10	9.851	164	476379	20.00	ng	0.00
64) Phenanthrene-d10	11.339	188	922550	20.00	ng	0.00
76) Chrysene-d12	13.986	240	657101	20.00	ng	0.00
86) Perylene-d12	15.433	264	499807	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1570325	88.31	ng	0.00
7) Phenol-d6	6.451	99	2179915	88.78	ng	0.00
23) Nitrobenzene-d5	7.381	82	2078805	91.97	ng	0.00
42) 2,4,6-Tribromophenol	10.645	330	573689	91.55	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	2956343	92.28	ng	0.00
79) Terphenyl-d14	12.927	244	3266911	90.99	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.516	88	435644	45.53	ng	97
3) Pyridine	3.251	79	1103190	46.83	ng	91
4) n-Nitrosodimethylamine	3.216	42	609306	46.97	ng	94
6) Aniline	6.475	93	1277770	44.77	ng	96
8) 2-Chlorophenol	6.598	128	868510	45.03	ng	96
9) Benzaldehyde	6.357	77	504354	42.71	ng	96
10) Phenol	6.463	94	1166607	42.41	ng	86
11) bis(2-Chloroethyl)ether	6.551	93	867598	44.92	ng	98
12) 1,3-Dichlorobenzene	6.751	146	859505	43.72	ng	99
13) 1,4-Dichlorobenzene	6.828	146	871983	44.14	ng	98
14) 1,2-Dichlorobenzene	6.986	146	795561	41.40	ng	100
15) Benzyl Alcohol	6.957	79	895427	45.98	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.086	45	1850950	45.07	ng	99
17) 2-Methylphenol	7.069	107	744824	44.92	ng	98
18) Hexachloroethane	7.322	117	352000	44.36	ng	93
19) n-Nitroso-di-n-propyla...	7.233	70	685557	44.13	ng	96
20) 3+4-Methylphenols	7.228	107	871134	45.33	ng	# 85
22) Acetophenone	7.222	105	1242258	44.63	ng	92
24) Nitrobenzene	7.398	77	1087707	46.15	ng	96
25) Isophorone	7.639	82	1937458	45.20	ng	99
26) 2-Nitrophenol	7.710	139	456688	46.77	ng	98
27) 2,4-Dimethylphenol	7.751	122	687260	45.15	ng	98
28) bis(2-Chloroethoxy)met...	7.851	93	985674	45.12	ng	98
29) 2,4-Dichlorophenol	7.957	162	671937	45.26	ng	99
30) 1,2,4-Trichlorobenzene	8.039	180	704933	45.13	ng	98
31) Naphthalene	8.122	128	2346349	44.33	ng	100
32) Benzoic acid	7.892	122	472758	49.86	ng	93
33) 4-Chloroaniline	8.169	127	1007594	45.20	ng	98
34) Hexachlorobutadiene	8.239	225	429582	44.88	ng	99
35) Caprolactam	8.551	113	244186m	46.47	ng	
36) 4-Chloro-3-methylphenol	8.657	107	861048	46.06	ng	94
37) 2-Methylnaphthalene	8.810	142	1524440	44.16	ng	100
38) 1-Methylnaphthalene	8.910	142	1427534	43.72	ng	99
40) 1,2,4,5-Tetrachloroben...	8.975	216	660817	45.56	ng	98
41) Hexachlorocyclopentadiene	8.963	237	298131	53.09	ng	98
43) 2,4,6-Trichlorophenol	9.086	196	467259	46.02	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.133	196	508987	46.99	ng	97
46) 1,1'-Biphenyl	9.275	154	1803124	44.91	ng	99
47) 2-Chloronaphthalene	9.298	162	1383587	45.38	ng	98
48) 2-Nitroaniline	9.392	65	641540	47.91	ng	91
49) Acenaphthylene	9.716	152	2227670	44.79	ng	99
50) Dimethylphthalate	9.580	163	1579754	45.28	ng	100
51) 2,6-Dinitrotoluene	9.639	165	377095	45.79	ng	96
52) Acenaphthene	9.886	154	1346788	44.78	ng	98
53) 3-Nitroaniline	9.810	138	443350	46.25	ng	99
54) 2,4-Dinitrophenol	9.916	184	223710	53.80	ng	90
55) Dibenzofuran	10.063	168	2013394	44.38	ng	99
56) 4-Nitrophenol	9.974	139	338510	47.23	ng	98
57) 2,4-Dinitrotoluene	10.045	165	513091	46.30	ng	91
58) Fluorene	10.404	166	1544043	44.15	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.180	232	394345	45.28	ng	99
60) Diethylphthalate	10.280	149	1683855	45.05	ng	98
61) 4-Chlorophenyl-phenyle...	10.392	204	711337	43.90	ng	97
62) 4-Nitroaniline	10.427	138	444174	46.42	ng	95
63) Azobenzene	10.557	77	1911779	44.85	ng	99
65) 4,6-Dinitro-2-methylph...	10.457	198	294503	47.73	ng	# 76
66) n-Nitrosodiphenylamine	10.516	169	1389148	44.11	ng	99
67) 4-Bromophenyl-phenylether	10.886	248	453991	43.88	ng	94
68) Hexachlorobenzene	10.951	284	527856	44.81	ng	98
69) Atrazine	11.045	200	432738	44.89	ng	98
70) Pentachlorophenol	11.145	266	294387	49.82	ng	95
71) Phenanthrene	11.363	178	2236900	43.90	ng	99
72) Anthracene	11.416	178	2224461	44.01	ng	99
73) Carbazole	11.574	167	2244906	43.70	ng	98
74) Di-n-butylphthalate	11.904	149	2536512	43.66	ng	99
75) Fluoranthene	12.557	202	2442490	43.96	ng	98
77) Benzidine	12.674	184	856963	44.12	ng	97
78) Pyrene	12.786	202	2470031	45.98	ng	98
80) Butylbenzylphthalate	13.404	149	1095642	47.27	ng	98
81) Benzo(a)anthracene	13.974	228	1920737	45.83	ng	98
82) 3,3'-Dichlorobenzidine	13.933	252	581238	45.40	ng	# 97
83) Chrysene	14.015	228	1922878	45.65	ng	99
84) Bis(2-ethylhexyl)phtha...	13.962	149	1328896	46.23	ng	97
85) Di-n-octyl phthalate	14.580	149	2157240	46.67	ng	99
87) Indeno(1,2,3-cd)pyrene	16.886	276	1433225	46.93	ng	99
88) Benzo(b)fluoranthene	15.015	252	1703533	48.33	ng	# 98
89) Benzo(k)fluoranthene	15.045	252	1486793	44.37	ng	98
90) Benzo(a)pyrene	15.380	252	1402766	46.20	ng	98
91) Dibenzo(a,h)anthracene	16.903	278	1207259	47.05	ng	98
92) Benzo(g,h,i)perylene	17.327	276	1230389	48.10	ng	97

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

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