

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081821\
 Data File : BP006753.D
 Acq On : 18 Aug 2021 20:00
 Operator : CG/JU
 Sample : PB138494BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138494BS

Manual Integrations
 APPROVED

mohammad
 8/19/2021 1:27:44 PM

Quant Time: Aug 19 04:50:05 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 18:17:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	168451	20.000	ng	0.00
21) Naphthalene-d8	10.481	136	705522	20.000	ng	#-0.01
39) Acenaphthene-d10	14.345	164	388269	20.000	ng	0.00
64) Phenanthrene-d10	17.104	188	750907	20.000	ng	# 0.00
76) Chrysene-d12	21.210	240	699667	20.000	ng	# 0.00
86) Perylene-d12	23.521	264	742838	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	1598591	145.760	ng	0.00
7) Phenol-d6	6.893	99	1978219	126.978	ng	0.00
23) Nitrobenzene-d5	8.863	82	1344184	87.941	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	599974	147.856	ng	0.00
45) 2-Fluorobiphenyl	12.963	172	2349662	90.540	ng	0.00
79) Terphenyl-d14	19.680	244	3326492	95.268	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.240	88	167510	35.083	ng	96
3) Pyridine	3.634	79	466987	33.276	ng	97
4) n-Nitrosodimethylamine	3.552	42	213588	40.324	ng	# 77
6) Aniline	7.040	93	768183	45.393	ng	96
8) 2-Chlorophenol	7.269	128	585728	49.360	ng	94
9) Benzaldehyde	6.852	77	409130	43.632	ng	93
10) Phenol	6.916	94	733299	46.941	ng	99
11) bis(2-Chloroethyl)ether	7.140	93	544686	45.919	ng	94
12) 1,3-Dichlorobenzene	7.587	146	583361	47.169	ng	97
13) 1,4-Dichlorobenzene	7.734	146	595078	47.410	ng	97
14) 1,2-Dichlorobenzene	8.046	146	571377	47.440	ng	96
15) Benzyl Alcohol	7.952	79	556705	47.650	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.234	45	608142	43.084	ng	95
17) 2-Methylphenol	8.152	107	488136	49.485	ng	99
18) Hexachloroethane	8.763	117	240594	48.428	ng	89
19) n-Nitroso-di-n-propyla...	8.516	70	458691	43.885	ng	96
20) 3+4-Methylphenols	8.481	107	660652	49.205	ng	97
22) Acetophenone	8.528	105	889166	47.524	ng	# 98
24) Nitrobenzene	8.905	77	703379	44.268	ng	92
25) Isophorone	9.434	82	1192588	43.440	ng	96
26) 2-Nitrophenol	9.610	139	298549	51.630	ng	93
27) 2,4-Dimethylphenol	9.681	122	521872	53.896	ng	94
28) bis(2-Chloroethoxy)met...	9.916	93	713866	48.597	ng	98
29) 2,4-Dichlorophenol	10.140	162	486784	49.655	ng	94
30) 1,2,4-Trichlorobenzene	10.352	180	493910	46.886	ng	97
31) Naphthalene	10.534	128	1730117	45.636	ng	99
32) Benzoic acid	9.846	122	377100	50.342	ng	90
33) 4-Chloroaniline	10.657	127	477878	31.150	ng	96
34) Hexachlorobutadiene	10.816	225	293146	47.755	ng	98
35) Caprolactam	11.469	113	180263m	50.995	ng	
36) 4-Chloro-3-methylphenol	11.793	107	606518	48.695	ng	89
37) 2-Methylnaphthalene	12.151	142	1169871	45.539	ng	100
38) 1-Methylnaphthalene	12.375	142	1082801	44.354	ng	99
40) 1,2,4,5-Tetrachloroben...	12.522	216	502500	49.448	ng	# 94
41) Hexachlorocyclopentadiene	12.498	237	695704	129.080	ng	99
43) 2,4,6-Trichlorophenol	12.769	196	357473	51.518	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.845	196	388984	50.644	ng	# 87
46) 1,1'-Biphenyl	13.175	154	1367892	46.545	ng	97
47) 2-Chloronaphthalene	13.216	162	1065309	47.072	ng	95
48) 2-Nitroaniline	13.428	65	400295	50.120	ng	# 85
49) Acenaphthylene	14.063	152	1774729	48.633	ng	99
50) Dimethylphthalate	13.816	163	1364800	48.290	ng	99
51) 2,6-Dinitrotoluene	13.934	165	302844	48.020	ng	# 80
52) Acenaphthene	14.410	154	1018691	45.868	ng	98
53) 3-Nitroaniline	14.263	138	258557	36.981	ng	83
54) 2,4-Dinitrophenol	14.475	184	338861	102.748	ng	# 78
55) Dibenzofuran	14.745	168	1578763	45.193	ng	93
56) 4-Nitrophenol	14.587	139	603979	99.489	ng	# 87
57) 2,4-Dinitrotoluene	14.728	165	422214	49.920	ng	# 75
58) Fluorene	15.398	166	1297837	46.481	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.975	232	317049	49.257	ng	# 69
60) Diethylphthalate	15.187	149	1455153	49.020	ng	98
61) 4-Chlorophenyl-phenyle...	15.398	204	598657	47.344	ng	# 89
62) 4-Nitroaniline	15.434	138	351238	47.343	ng	93
63) Azobenzene	15.692	77	1600173	46.235	ng	93
65) 4,6-Dinitro-2-methylph...	15.492	198	204360	44.540	ng	75
66) n-Nitrosodiphenylamine	15.616	169	1114570	47.223	ng	100
67) 4-Bromophenyl-phenylether	16.298	248	355251	49.307	ng	# 88
68) Hexachlorobenzene	16.398	284	401286	48.943	ng	# 85
69) Atrazine	16.586	200	399956	54.305	ng	96
70) Pentachlorophenol	16.751	266	542838	104.350	ng	99
71) Phenanthrene	17.145	178	1984115	46.844	ng	99
72) Anthracene	17.239	178	2031427	49.635	ng	99
73) Carbazole	17.516	167	1903835	45.605	ng	98
74) Di-n-butylphthalate	18.081	149	2491382	52.382	ng	99
75) Fluoranthene	19.122	202	2231478	46.939	ng	95
77) Benzidine	19.316	184	1572699	89.827	ng	98
78) Pyrene	19.475	202	2296056	49.130	ng	99
80) Butylbenzylphthalate	20.369	149	1137101	55.332	ng	87
81) Benzo(a)anthracene	21.192	228	2164549	47.338	ng	99
82) 3,3'-Dichlorobenzidine	21.133	252	606099	50.490	ng	# 97
83) Chrysene	21.245	228	2106511	47.520	ng	99
84) Bis(2-ethylhexyl)phtha...	21.133	149	1711422	55.418	ng	# 96
85) Di-n-octyl phthalate	22.033	149	2966120	58.272	ng	# 96
87) Indeno(1,2,3-cd)pyrene	25.910	276	2694331	50.021	ng	# 91
88) Benzo(b)fluoranthene	22.821	252	2330395	48.270	ng	# 97
89) Benzo(k)fluoranthene	22.869	252	2158744	46.571	ng	# 97
90) Benzo(a)pyrene	23.421	252	2212457	51.092	ng	# 96
91) Dibenzo(a,h)anthracene	25.927	278	2299030	49.366	ng	# 94
92) Benzo(g,h,i)perylene	26.645	276	2244501	50.294	ng	# 91

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

