

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091621\
 Data File : BP007000.D
 Acq On : 16 Sep 2021 17:46
 Operator : CG/JU
 Sample : PB139068BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139068BS

Manual Integrations
 APPROVED

mohammad
 9/17/2021 11:29:47 AM

Quant Time: Sep 16 18:13:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 16 15:16:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	363106	20.000	ng	-0.01
21) Naphthalene-d8	10.716	136	1473106	20.000	ng	# 0.00
39) Acenaphthene-d10	14.539	164	944605	20.000	ng	0.00
64) Phenanthrene-d10	17.286	188	1953393	20.000	ng	# 0.00
76) Chrysene-d12	21.368	240	1918775	20.000	ng	#-0.01
86) Perylene-d12	23.780	264	1924862	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	2837528	127.607	ng	0.00
7) Phenol-d6	7.081	99	3511777	115.706	ng	0.00
23) Nitrobenzene-d5	9.075	82	1972776	75.216	ng	0.00
42) 2,4,6-Tribromophenol	16.028	330	1381242	121.589	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	4799504	70.143	ng	0.00
79) Terphenyl-d14	19.839	244	7415817	74.152	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.375	88	323596	33.234	ng	# 87
3) Pyridine	3.775	79	835961	33.391	ng	# 85
4) n-Nitrosodimethylamine	3.687	42	397677	43.240	ng	# 81
6) Aniline	7.240	93	1379905	42.194	ng	99
8) 2-Chlorophenol	7.475	128	1033755	41.478	ng	97
9) Benzaldehyde	7.052	77	619510m	44.778	ng	
10) Phenol	7.105	94	1303952	42.619	ng	91
11) bis(2-Chloroethyl)ether	7.340	93	964038	40.877	ng	95
12) 1,3-Dichlorobenzene	7.805	146	1090796	39.062	ng	96
13) 1,4-Dichlorobenzene	7.946	146	1116091	39.420	ng	99
14) 1,2-Dichlorobenzene	8.263	146	1068530	39.413	ng	98
15) Benzyl Alcohol	8.152	79	835603	42.686	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.446	45	1142801	41.451	ng	88
17) 2-Methylphenol	8.352	107	853785	43.443	ng	94
18) Hexachloroethane	8.993	117	384780	40.921	ng	98
19) n-Nitroso-di-n-propyla...	8.722	70	671373	40.291	ng	99
20) 3+4-Methylphenols	8.681	107	1155353	43.432	ng	94
22) Acetophenone	8.740	105	1460929	39.322	ng	99
24) Nitrobenzene	9.116	77	1026744	37.523	ng	97
25) Isophorone	9.646	82	1823666	37.614	ng	98
26) 2-Nitrophenol	9.828	139	500995	42.230	ng	# 89
27) 2,4-Dimethylphenol	9.887	122	961673	46.066	ng	96
28) bis(2-Chloroethoxy)met...	10.128	93	1228712	41.769	ng	99
29) 2,4-Dichlorophenol	10.357	162	969393	40.646	ng	99
30) 1,2,4-Trichlorobenzene	10.575	180	1014419	37.405	ng	98
31) Naphthalene	10.763	128	3008943	36.864	ng	100
32) Benzoic acid	10.028	122	679838	44.284	ng	95
33) 4-Chloroaniline	10.869	127	924660	29.176	ng	99
34) Hexachlorobutadiene	11.051	225	579739	36.299	ng	99
35) Caprolactam	11.669	113	309194m	47.115	ng	
36) 4-Chloro-3-methylphenol	11.993	107	1014289	42.158	ng	98
37) 2-Methylnaphthalene	12.369	142	2193551	37.907	ng	96
38) 1-Methylnaphthalene	12.587	142	2009881	36.783	ng	100
40) 1,2,4,5-Tetrachloroben...	12.740	216	1119181	37.231	ng	99
41) Hexachlorocyclopentadiene	12.722	237	1445318	96.790	ng	98
43) 2,4,6-Trichlorophenol	12.975	196	791510	40.731	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.040	196	865181	40.409	ng	98
46) 1,1'-Biphenyl	13.375	154	2666838	36.194	ng	99
47) 2-Chloronaphthalene	13.416	162	2140775	36.875	ng	98
48) 2-Nitroaniline	13.616	65	607041	45.634	ng	99
49) Acenaphthylene	14.263	152	3457920	39.306	ng	100
50) Dimethylphthalate	14.004	163	2724326	39.515	ng	100
51) 2,6-Dinitrotoluene	14.116	165	595233	39.671	ng	94
52) Acenaphthene	14.604	154	2057095	36.943	ng	99
53) 3-Nitroaniline	14.445	138	524411	34.964	ng	99
54) 2,4-Dinitrophenol	14.651	184	700862	88.090	ng	90
55) Dibenzofuran	14.939	168	3285548	36.429	ng	95
56) 4-Nitrophenol	14.751	139	1142732	89.772	ng	# 81
57) 2,4-Dinitrotoluene	14.898	165	816273	42.505	ng	97
58) Fluorene	15.586	166	2660733	37.798	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.163	232	737411m	40.042	ng	
60) Diethylphthalate	15.363	149	2674589	40.337	ng	98
61) 4-Chlorophenyl-phenyle...	15.581	204	1348440	38.192	ng	92
62) 4-Nitroaniline	15.604	138	637390	42.275	ng	# 81
63) Azobenzene	15.875	77	2386421	38.863	ng	94
65) 4,6-Dinitro-2-methylph...	15.663	198	435737	36.784	ng	99
66) n-Nitrosodiphenylamine	15.798	169	2373296	38.176	ng	99
67) 4-Bromophenyl-phenylether	16.475	248	831221	38.495	ng	96
68) Hexachlorobenzene	16.592	284	930627	38.337	ng	98
69) Atrazine	16.751	200	856367	46.327	ng	97
70) Pentachlorophenol	16.933	266	1279136	87.744	ng	100
71) Phenanthrene	17.328	178	4228261	37.312	ng	99
72) Anthracene	17.422	178	4318077	39.826	ng	99
73) Carbazole	17.692	167	3941230	37.262	ng	98
74) Di-n-butylphthalate	18.245	149	4691765	43.372	ng	99
75) Fluoranthene	19.292	202	4944802	37.884	ng	97
77) Benzidine	19.469	184	2669792	77.864	ng	99
78) Pyrene	19.645	202	5106246	38.404	ng	99
80) Butylbenzylphthalate	20.516	149	2087685	44.849	ng	98
81) Benzo(a)anthracene	21.351	228	4902664	37.602	ng	98
82) 3,3'-Dichlorobenzidine	21.286	252	1643580	44.332	ng	98
83) Chrysene	21.404	228	4827029	37.728	ng	98
84) Bis(2-ethylhexyl)phtha...	21.280	149	3020798	44.148	ng	97
85) Di-n-octyl phthalate	22.204	149	5170703	46.235	ng	99
87) Indeno(1,2,3-cd)pyrene	26.315	276	5927962	40.580	ng	# 94
88) Benzo(b)fluoranthene	23.045	252	5240156	40.081	ng	# 98
89) Benzo(k)fluoranthene	23.092	252	4971442	39.068	ng	# 98
90) Benzo(a)pyrene	23.674	252	5005358	42.691	ng	# 97
91) Dibenzo(a,h)anthracene	26.333	278	5064850	40.080	ng	# 95
92) Benzo(g,h,i)perylene	27.092	276	4993334	40.957	ng	# 94

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

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