

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081921\
 Data File : BP006776.D
 Acq On : 19 Aug 2021 12:05
 Operator : CG/JU
 Sample : PB138536BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138536BS

Manual Integrations
 APPROVED

mohammad
 8/20/2021 2:24:25 PM

Quant Time: Aug 19 15:15:27 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	7.693	152	144396	20.000	ng	-0.01
21) Naphthalene-d8	10.481	136	625652	20.000	ng	#-0.01
39) Acenaphthene-d10	14.340	164	349614	20.000	ng	0.00
64) Phenanthrene-d10	17.098	188	688296	20.000	ng	# 0.00
76) Chrysene-d12	21.204	240	669237	20.000	ng	# 0.00
86) Perylene-d12	23.516	264	689655	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	1399630	148.879	ng	0.00
7) Phenol-d6	6.887	99	1726677	129.296	ng	0.00
23) Nitrobenzene-d5	8.858	82	1185750	87.479	ng	0.00
42) 2,4,6-Tribromophenol	15.840	330	524587	143.571	ng	0.00
45) 2-Fluorobiphenyl	12.963	172	2075497	88.818	ng	0.00
79) Terphenyl-d14	19.681	244	3022654	90.502	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.240	88	161838	39.541	ng	97
3) Pyridine	3.634	79	417125	34.675	ng	98
4) n-Nitrosodimethylamine	3.552	42	198566	43.733	ng	# 75
6) Aniline	7.034	93	683786	47.137	ng	96
8) 2-Chlorophenol	7.264	128	503562	49.505	ng	92
9) Benzaldehyde	6.846	77	378927	47.144	ng	93
10) Phenol	6.911	94	637976	47.642	ng	96
11) bis(2-Chloroethyl)ether	7.134	93	475524	46.767	ng	93
12) 1,3-Dichlorobenzene	7.581	146	502616	47.410	ng	96
13) 1,4-Dichlorobenzene	7.728	146	513026	47.682	ng	97
14) 1,2-Dichlorobenzene	8.046	146	493246	47.775	ng	97
15) Benzyl Alcohol	7.946	79	479230	47.852	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.228	45	531031	43.888	ng	94
17) 2-Methylphenol	8.152	107	421329	49.828	ng	97
18) Hexachloroethane	8.764	117	212173	49.822	ng	90
19) n-Nitroso-di-n-propyla...	8.511	70	401129	44.771	ng	# 96
20) 3+4-Methylphenols	8.481	107	570771	49.593	ng	97
22) Acetophenone	8.522	105	774939	46.706	ng	# 97
24) Nitrobenzene	8.899	77	624225	44.302	ng	90
25) Isophorone	9.428	82	1033151	42.436	ng	96
26) 2-Nitrophenol	9.605	139	249334	48.624	ng	92
27) 2,4-Dimethylphenol	9.675	122	455162	53.007	ng	96
28) bis(2-Chloroethoxy)met...	9.911	93	628983	48.285	ng	98
29) 2,4-Dichlorophenol	10.134	162	418184	48.103	ng	94
30) 1,2,4-Trichlorobenzene	10.346	180	429807	46.010	ng	97
31) Naphthalene	10.528	128	1513875	45.030	ng	99
32) Benzoic acid	9.840	122	300351	45.214	ng	92
33) 4-Chloroaniline	10.652	127	469996	34.547	ng	# 95
34) Hexachlorobutadiene	10.810	225	256196	47.063	ng	99
35) Caprolactam	11.463	113	147740m	47.130	ng	
36) 4-Chloro-3-methylphenol	11.787	107	525164	47.546	ng	89
37) 2-Methylnaphthalene	12.146	142	1030249	45.224	ng	98
38) 1-Methylnaphthalene	12.369	142	948661	43.820	ng	98
40) 1,2,4,5-Tetrachloroben...	12.522	216	440132	48.100	ng	# 94
41) Hexachlorocyclopentadiene	12.493	237	589348	121.437	ng	100
43) 2,4,6-Trichlorophenol	12.769	196	307298	49.183	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.840	196	333789	48.263	ng	# 86
46) 1,1'-Biphenyl	13.169	154	1208215	45.658	ng	97
47) 2-Chloronaphthalene	13.210	162	937978	46.028	ng	94
48) 2-Nitroaniline	13.428	65	350667	48.760	ng	85
49) Acenaphthylene	14.057	152	1564732	47.619	ng	100
50) Dimethylphthalate	13.816	163	1192822	46.871	ng	99
51) 2,6-Dinitrotoluene	13.934	165	258709	45.557	ng	# 75
52) Acenaphthene	14.404	154	901606	45.085	ng	96
53) 3-Nitroaniline	14.263	138	243762	38.720	ng	82
54) 2,4-Dinitrophenol	14.475	184	287437	96.792	ng	# 79
55) Dibenzofuran	14.746	168	1412902	44.917	ng	94
56) 4-Nitrophenol	14.587	139	531862	97.296	ng	89
57) 2,4-Dinitrotoluene	14.722	165	369665	48.540	ng	# 71
58) Fluorene	15.393	166	1155038	45.940	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.975	232	274784	47.411	ng	# 67
60) Diethylphthalate	15.187	149	1291924	48.333	ng	97
61) 4-Chlorophenyl-phenyle...	15.393	204	529997	46.548	ng	# 87
62) 4-Nitroaniline	15.434	138	287897m	43.096	ng	
63) Azobenzene	15.687	77	1468969	47.137	ng	92
65) 4,6-Dinitro-2-methylph...	15.487	198	175265	41.674	ng	69
66) n-Nitrosodiphenylamine	15.616	169	990490	45.783	ng	99
67) 4-Bromophenyl-phenylether	16.293	248	313239	47.431	ng	# 86
68) Hexachlorobenzene	16.398	284	353043	46.976	ng	# 86
69) Atrazine	16.581	200	351099	52.008	ng	96
70) Pentachlorophenol	16.751	266	482652	101.220	ng	99
71) Phenanthrene	17.140	178	1791933	46.155	ng	99
72) Anthracene	17.234	178	1818855	48.483	ng	100
73) Carbazole	17.510	167	1716628	44.861	ng	98
74) Di-n-butylphthalate	18.081	149	2196323	50.379	ng	99
75) Fluoranthene	19.122	202	2007069	46.059	ng	95
77) Benzidine	19.316	184	1458147	87.072	ng	99
78) Pyrene	19.475	202	2061256	46.112	ng	98
80) Butylbenzylphthalate	20.369	149	1000062	50.877	ng	90
81) Benzo(a)anthracene	21.192	228	1965986	44.950	ng	99
82) 3,3'-Dichlorobenzidine	21.133	252	591817	51.542	ng	# 97
83) Chrysene	21.239	228	1931018	45.542	ng	99
84) Bis(2-ethylhexyl)phtha...	21.133	149	1497718	50.703	ng	# 97
85) Di-n-octyl phthalate	22.033	149	2584246	53.079	ng	# 95
87) Indeno(1,2,3-cd)pyrene	25.904	276	2436671	48.726	ng	# 90
88) Benzo(b)fluoranthene	22.816	252	2092622	46.687	ng	# 96
89) Benzo(k)fluoranthene	22.863	252	1995694	46.374	ng	# 96
90) Benzo(a)pyrene	23.416	252	1998192	49.703	ng	# 95
91) Dibenzo(a,h)anthracene	25.921	278	2084809	48.219	ng	# 95
92) Benzo(g,h,i)perylene	26.639	276	2028232	48.952	ng	# 91

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

