

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080521\
 Data File : BP006501.D
 Acq On : 05 Aug 2021 11:23
 Operator : CG/JU
 Sample : PB138179BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138179BS

Manual Integrations
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Quant Time: Aug 05 11:57:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 15:41:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	160892	20.000	ng	0.00
21) Naphthalene-d8	10.540	136	694736	20.000	ng	# 0.00
39) Acenaphthene-d10	14.386	164	381654	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	718931	20.000	ng	# 0.00
76) Chrysene-d12	21.239	240	641562	20.000	ng	# 0.00
86) Perylene-d12	23.562	264	638771	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	1517620	144.878	ng	0.00
7) Phenol-d6	6.934	99	2021320	135.840	ng	0.00
23) Nitrobenzene-d5	8.910	82	1139269	75.692	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	486943	122.080	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	1922199	75.352	ng	0.00
79) Terphenyl-d14	19.716	244	2239266	69.939	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.305	88	226070	49.572	ng	90
3) Pyridine	3.693	79	572443	42.707	ng	95
4) n-Nitrosodimethylamine	3.605	42	273246	54.011	ng	# 92
6) Aniline	7.093	93	768643	47.554	ng	94
8) 2-Chlorophenol	7.322	128	596939	52.668	ng	88
9) Benzaldehyde	6.904	77	338120	37.754	ng	89
10) Phenol	6.963	94	782324	52.432	ng	97
11) bis(2-Chloroethyl)ether	7.193	93	578303	51.044	ng	89
12) 1,3-Dichlorobenzene	7.646	146	590255	49.968	ng	# 95
13) 1,4-Dichlorobenzene	7.793	146	601740	50.193	ng	97
14) 1,2-Dichlorobenzene	8.104	146	578121	50.255	ng	96
15) Benzyl Alcohol	8.004	79	580580	52.028	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.293	45	694221	51.493	ng	92
17) 2-Methylphenol	8.204	107	506161	53.723	ng	99
18) Hexachloroethane	8.828	117	249096	52.495	ng	# 86
19) n-Nitroso-di-n-propyla...	8.569	70	488425	48.925	ng	# 95
20) 3+4-Methylphenols	8.528	107	672412	52.434	ng	96
22) Acetophenone	8.581	105	894154	48.532	ng	# 96
24) Nitrobenzene	8.957	77	741715	47.406	ng	90
25) Isophorone	9.481	82	1200095	44.392	ng	94
26) 2-Nitrophenol	9.663	139	274133	48.144	ng	91
27) 2,4-Dimethylphenol	9.728	122	541212	56.761	ng	96
28) bis(2-Chloroethoxy)met...	9.969	93	755440	52.226	ng	# 98
29) 2,4-Dichlorophenol	10.193	162	476001	49.309	ng	94
30) 1,2,4-Trichlorobenzene	10.404	180	490201	47.257	ng	95
31) Naphthalene	10.592	128	1756829	47.060	ng	99
32) Benzoic acid	9.857	122	328190	44.493	ng	87
33) 4-Chloroaniline	10.704	127	302804	20.044	ng	# 93
34) Hexachlorobutadiene	10.875	225	278968	46.150	ng	98
35) Caprolactam	11.498	113	158535m	45.545	ng	
36) 4-Chloro-3-methylphenol	11.834	107	589286	48.046	ng	88
37) 2-Methylnaphthalene	12.204	142	1201222	47.486	ng	99
38) 1-Methylnaphthalene	12.422	142	1100637	45.784	ng	98
40) 1,2,4,5-Tetrachloroben...	12.575	216	488919	48.946	ng	# 94
41) Hexachlorocyclopentadiene	12.551	237	719569	135.822	ng	98
43) 2,4,6-Trichlorophenol	12.816	196	331671	48.628	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.886	196	364202	48.240	ng	# 89
46) 1,1'-Biphenyl	13.222	154	1398264	48.403	ng	96
47) 2-Chloronaphthalene	13.263	162	1082949	48.681	ng	94
48) 2-Nitroaniline	13.469	65	399080	50.834	ng	# 84
49) Acenaphthylene	14.110	152	1770746	49.365	ng	99
50) Dimethylphthalate	13.857	163	1299392	46.772	ng	99
51) 2,6-Dinitrotoluene	13.975	165	281939	45.480	ng	# 77
52) Acenaphthene	14.451	154	1040521	47.663	ng	98
53) 3-Nitroaniline	14.298	138	208431	30.329	ng	82
54) 2,4-Dinitrophenol	14.504	184	278591	85.937	ng	# 73
55) Dibenzofuran	14.786	168	1592679	46.382	ng	93
56) 4-Nitrophenol	14.610	139	556697	93.290	ng	# 85
57) 2,4-Dinitrotoluene	14.757	165	389255	46.821	ng	# 76
58) Fluorene	15.439	166	1290258	47.010	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.010	232	288140	45.541	ng	# 67
60) Diethylphthalate	15.228	149	1368802	46.910	ng	98
61) 4-Chlorophenyl-phenyle...	15.439	204	582939	46.900	ng	# 87
62) 4-Nitroaniline	15.463	138	321072	44.027	ng	91
63) Azobenzene	15.733	77	1643730	48.317	ng	92
65) 4,6-Dinitro-2-methylph...	15.516	198	169931	38.684	ng	# 66
66) n-Nitrosodiphenylamine	15.657	169	1092773	48.359	ng	99
67) 4-Bromophenyl-phenylether	16.333	248	329667	47.791	ng	# 85
68) Hexachlorobenzene	16.439	284	373734	47.610	ng	# 85
69) Atrazine	16.616	200	347507	49.282	ng	95
70) Pentachlorophenol	16.786	266	485176	97.413	ng	98
71) Phenanthrene	17.180	178	1935897	47.739	ng	99
72) Anthracene	17.274	178	1942657	49.577	ng	100
73) Carbazole	17.551	167	1798297	44.993	ng	98
74) Di-n-butylphthalate	18.116	149	2192637	48.152	ng	99
75) Fluoranthene	19.157	202	2022201	44.429	ng	94
77) Benzidine	19.345	184	961230	59.875	ng	99
78) Pyrene	19.510	202	2105477	49.133	ng	99
80) Butylbenzylphthalate	20.398	149	926731	49.180	ng	# 84
81) Benzo(a)anthracene	21.221	228	1911694	45.594	ng	99
82) 3,3'-Dichlorobenzidine	21.163	252	496306	45.089	ng	# 97
83) Chrysene	21.274	228	1883676	46.341	ng	99
84) Bis(2-ethylhexyl)phtha...	21.163	149	1361286	48.072	ng	# 95
85) Di-n-octyl phthalate	22.068	149	2241393	48.022	ng	# 93
87) Indeno(1,2,3-cd)pyrene	25.974	276	2297028	49.593	ng	# 88
88) Benzo(b)fluoranthene	22.862	252	1912090	46.058	ng	# 96
89) Benzo(k)fluoranthene	22.904	252	1923288	48.252	ng	# 95
90) Benzo(a)pyrene	23.462	252	1860617	49.967	ng	# 94
91) Dibenzo(a,h)anthracene	25.998	278	1975954	49.341	ng	# 93
92) Benzo(g,h,i)perylene	26.715	276	1945441	50.694	ng	# 89

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

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