

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
 Data File : BF124131.D
 Acq On : 4 Jun 2021 10:22
 Operator : JU/CG
 Sample : PB136792BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB136792BS

Manual Integrations
 APPROVED

mohammad
 6/7/2021 12:24:08 PM

Quant Time: Jun 04 10:55:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 04 09:56:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	52018	20.00	ng	0.00
21) Naphthalene-d8	8.151	136	198073	20.00	ng	0.00
39) Acenaphthene-d10	9.910	164	116428	20.00	ng	0.00
64) Phenanthrene-d10	11.392	188	205942	20.00	ng	0.00
76) Chrysene-d12	14.039	240	133114	20.00	ng	0.00
86) Perylene-d12	15.510	264	98810	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	401187	116.10	ng	0.02
7) Phenol-d6	6.516	99	508907	109.56	ng	0.00
23) Nitrobenzene-d5	7.434	82	382166	76.52	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	187709	113.80	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	717127	77.51	ng	0.00
79) Terphenyl-d14	12.980	244	779193	80.37	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.746	88	51913	34.33	ng	95
3) Pyridine	3.493	79	125814	32.48	ng	91
4) n-Nitrosodimethylamine	3.457	42	94911	41.24	ng	95
6) Aniline	6.528	93	161985	30.53	ng	# 1
8) 2-Chlorophenol	6.657	128	159267	41.42	ng	94
9) Benzaldehyde	6.416	77	71765	34.76	ng	96
10) Phenol	6.528	94	200439	39.93	ng	# 69
11) bis(2-Chloroethyl)ether	6.604	93	153536	41.96	ng	89
12) 1,3-Dichlorobenzene	6.810	146	177225m	39.43	ng	
13) 1,4-Dichlorobenzene	6.887	146	183751	40.75	ng	94
14) 1,2-Dichlorobenzene	7.040	146	172684	40.04	ng	98
15) Benzyl Alcohol	7.016	79	157515	37.62	ng	92
16) 2,2'-oxybis(1-Chloropr...	7.145	45	231270	40.51	ng	88
17) 2-Methylphenol	7.128	107	128538	41.12	ng	88
18) Hexachloroethane	7.381	117	71066	40.02	ng	# 87
19) n-Nitroso-di-n-propyla...	7.287	70	128086	39.06	ng	88
20) 3+4-Methylphenols	7.281	107	166567	40.22	ng	# 86
22) Acetophenone	7.275	105	236144	41.82	ng	# 94
24) Nitrobenzene	7.451	77	190037	37.93	ng	96
25) Isophorone	7.692	82	327930	36.44	ng	99
26) 2-Nitrophenol	7.769	139	87114	38.93	ng	# 88
27) 2,4-Dimethylphenol	7.810	122	139102	44.08	ng	93
28) bis(2-Chloroethoxy)met...	7.898	93	186143	41.72	ng	98
29) 2,4-Dichlorophenol	8.016	162	143492	39.30	ng	89
30) 1,2,4-Trichlorobenzene	8.092	180	160366	39.19	ng	96
31) Naphthalene	8.175	128	446629	37.72	ng	99
32) Benzoic acid	7.957	122	74336	38.39	ng	85
33) 4-Chloroaniline	8.216	127	45599	10.35	ng	99
34) Hexachlorobutadiene	8.287	225	107667	37.25	ng	98
35) Caprolactam	8.604	113	41170m	41.04	ng	
36) 4-Chloro-3-methylphenol	8.710	107	152965	38.29	ng	87
37) 2-Methylnaphthalene	8.863	142	326481	39.31	ng	99
38) 1-Methylnaphthalene	8.963	142	299807	38.63	ng	97
40) 1,2,4,5-Tetrachloroben...	9.034	216	175106	42.45	ng	97
41) Hexachlorocyclopentadiene	9.022	237	217549	95.57	ng	94
43) 2,4,6-Trichlorophenol	9.145	196	108085	40.17	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	113953	39.46	ng	94
46) 1,1'-Biphenyl	9.328	154	414207	42.14	ng	97
47) 2-Chloronaphthalene	9.357	162	317261	39.66	ng	95
48) 2-Nitroaniline	9.451	65	110412	38.24	ng	98
49) Acenaphthylene	9.775	152	482477	41.49	ng	98
50) Dimethylphthalate	9.633	163	384982	39.97	ng	99
51) 2,6-Dinitrotoluene	9.692	165	85357	38.65	ng	97
52) Acenaphthene	9.945	154	288493	37.98	ng	92
53) 3-Nitroaniline	9.863	138	43381	20.99	ng	99
54) 2,4-Dinitrophenol	9.975	184	93450	82.64	ng	87
55) Dibenzofuran	10.116	168	459361	38.64	ng	99
56) 4-Nitrophenol	10.039	139	121925	82.53	ng	# 83
57) 2,4-Dinitrotoluene	10.098	165	114920	39.45	ng	# 94
58) Fluorene	10.457	166	361481	39.62	ng	95
59) 2,3,4,6-Tetrachlorophenol	10.233	232	96186	41.10	ng	# 90
60) Diethylphthalate	10.333	149	387789	39.89	ng	95
61) 4-Chlorophenyl-phenyle...	10.445	204	183458	39.32	ng	95
62) 4-Nitroaniline	10.486	138	77198	37.14	ng	# 79
63) Azobenzene	10.610	77	374032	36.91	ng	95
65) 4,6-Dinitro-2-methylph...	10.510	198	61754	37.31	ng	# 78
66) n-Nitrosodiphenylamine	10.569	169	316686	40.47	ng	98
67) 4-Bromophenyl-phenylether	10.939	248	115109	40.72	ng	96
68) Hexachlorobenzene	11.004	284	131780	41.16	ng	97
69) Atrazine	11.098	200	114097	51.51	ng	98
70) Pentachlorophenol	11.204	266	129392	76.89	ng	97
71) Phenanthrene	11.422	178	528283	41.06	ng	97
72) Anthracene	11.475	178	540431	41.71	ng	98
73) Carbazole	11.627	167	434801	38.51	ng	95
74) Di-n-butylphthalate	11.951	149	604172	41.64	ng	99
75) Fluoranthene	12.610	202	534091	39.55	ng	98
77) Benzidine	12.727	184	173207m	55.17	ng	
78) Pyrene	12.839	202	527234	41.25	ng	98
80) Butylbenzylphthalate	13.451	149	219995	42.43	ng	96
81) Benzo(a)anthracene	14.027	228	390146	40.23	ng	98
82) 3,3'-Dichlorobenzidine	13.986	252	70149	24.66	ng	98
83) Chrysene	14.063	228	370479	39.60	ng	99
84) Bis(2-ethylhexyl)phtha...	14.010	149	276735	44.87	ng	94
85) Di-n-octyl phthalate	14.621	149	369083	44.84	ng	# 92
87) Indeno(1,2,3-cd)pyrene	17.009	276	325287	40.32	ng	99
88) Benzo(b)fluoranthene	15.080	252	318393	41.48	ng	99
89) Benzo(k)fluoranthene	15.115	252	280597	38.62	ng	98
90) Benzo(a)pyrene	15.451	252	277806	41.67	ng	97
91) Dibenzo(a,h)anthracene	17.021	278	274804	39.93	ng	96
92) Benzo(g,h,i)perylene	17.457	276	272382	40.01	ng	96

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

