

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102120\
 Data File : BF122049.D
 Acq On : 21 Oct 2020 11:59
 Operator : JU/CG
 Sample : PB132505BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB132505BS

Manual Integrations
 APPROVED

mohammad
 10/22/2020 12:59:00 PM

Quant Time: Oct 21 12:29:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 12:04:42 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.728	152	120182	20.00	ng	0.00
21) Naphthalene-d8	8.010	136	462534	20.00	ng	0.00
39) Acenaphthene-d10	9.763	164	243140	20.00	ng	0.00
64) Phenanthrene-d10	11.245	188	458665	20.00	ng	0.00
76) Chrysene-d12	13.886	240	338673	20.00	ng	0.00
86) Perylene-d12	15.310	264	288636	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	928621	114.04	ng	0.02
7) Phenol-d6	6.387	99	1190223	113.55	ng	0.00
23) Nitrobenzene-d5	7.298	82	707097	78.40	ng	0.00
42) 2,4,6-Tribromophenol	10.557	330	377086	125.37	ng	0.00
45) 2-Fluorobiphenyl	9.086	172	1240345	75.06	ng	0.00
79) Terphenyl-d14	12.827	244	1405719	79.10	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.540	88	158105	44.15	ng	98
3) Pyridine	3.257	79	428434	45.50	ng	99
4) n-Nitrosodimethylamine	3.222	42	214698	47.17	ng	95
6) Aniline	6.398	93	389845	30.20	ng	# 1
8) 2-Chlorophenol	6.522	128	380949	46.29	ng	97
9) Benzaldehyde	6.281	77	180766	28.30	ng	100
10) Phenol	6.398	94	470978	43.54	ng	96
11) bis(2-Chloroethyl)ether	6.475	93	392566	46.94	ng	95
12) 1,3-Dichlorobenzene	6.669	146	414793	44.81	ng	98
13) 1,4-Dichlorobenzene	6.745	146	419570	45.14	ng	98
14) 1,2-Dichlorobenzene	6.898	146	383573	45.14	ng	98
15) Benzyl Alcohol	6.887	79	324288	47.83	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.010	45	574347	44.63	ng	70
17) 2-Methylphenol	6.998	107	318182	45.22	ng	# 89
18) Hexachloroethane	7.234	117	156556	46.63	ng	97
19) n-Nitroso-di-n-propyla...	7.151	70	281118	47.10	ng	100
20) 3+4-Methylphenols	7.157	107	398706	46.99	ng	# 68
22) Acetophenone	7.145	105	536113	45.05	ng	92
24) Nitrobenzene	7.322	77	442301	45.88	ng	97
25) Isophorone	7.557	82	790689	46.37	ng	99
26) 2-Nitrophenol	7.634	139	210651	47.45	ng	95
27) 2,4-Dimethylphenol	7.681	122	342541	45.26	ng	98
28) bis(2-Chloroethoxy)met...	7.763	93	492792	46.19	ng	99
29) 2,4-Dichlorophenol	7.881	162	328317	46.33	ng	99
30) 1,2,4-Trichlorobenzene	7.951	180	341105	45.27	ng	99
31) Naphthalene	8.034	128	1108984	44.76	ng	100
32) Benzoic acid	7.845	122	202786	47.06	ng	96
33) 4-Chloroaniline	8.092	127	233683	22.14	ng	97
34) Hexachlorobutadiene	8.145	225	207537	46.45	ng	99
35) Caprolactam	8.481	113	102496m	45.54	ng	
36) 4-Chloro-3-methylphenol	8.592	107	360918	47.42	ng	97
37) 2-Methylnaphthalene	8.722	142	725304	45.20	ng	98
38) 1-Methylnaphthalene	8.822	142	668487	44.98	ng	99
40) 1,2,4,5-Tetrachloroben...	8.892	216	348472	44.41	ng	100
41) Hexachlorocyclopentadiene	8.869	237	187408	75.84	ng	98
43) 2,4,6-Trichlorophenol	9.010	196	221602	45.77	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.057	196	232451	44.45	ng	98
46) 1,1'-Biphenyl	9.186	154	877203	44.37	ng	99
47) 2-Chloronaphthalene	9.210	162	689690	44.86	ng	99
48) 2-Nitroaniline	9.322	65	234121	46.17	ng	94
49) Acenaphthylene	9.628	152	1019155	43.16	ng	100
50) Dimethylphthalate	9.492	163	820880	46.17	ng	100
51) 2,6-Dinitrotoluene	9.563	165	178576	45.79	ng	97
52) Acenaphthene	9.798	154	620299	44.16	ng	98
53) 3-Nitroaniline	9.733	138	114730	25.87	ng	97
54) 2,4-Dinitrophenol	9.857	184	161993	81.24	ng	96
55) Dibenzofuran	9.969	168	894489	43.55	ng	97
56) 4-Nitrophenol	9.922	139	178890	74.14	ng	86
57) 2,4-Dinitrotoluene	9.975	165	220459	45.92	ng	97
58) Fluorene	10.310	166	738556	44.63	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.098	232	194786	48.15	ng	98
60) Diethylphthalate	10.192	149	810861	47.30	ng	99
61) 4-Chlorophenyl-phenyle...	10.304	204	363405	45.79	ng	99
62) 4-Nitroaniline	10.357	138	173972	39.26	ng	98
63) Azobenzene	10.463	77	826074	45.56	ng	99
65) 4,6-Dinitro-2-methylph...	10.386	198	129381	42.64	ng	99
66) n-Nitrosodiphenylamine	10.428	169	696045	44.05	ng	100
67) 4-Bromophenyl-phenylether	10.792	248	243764	46.00	ng	97
68) Hexachlorobenzene	10.863	284	272527	45.44	ng	93
69) Atrazine	10.957	200	192392	49.80	ng	98
70) Pentachlorophenol	11.063	266	206353	86.76	ng	99
71) Phenanthrene	11.275	178	1086259	43.19	ng	99
72) Anthracene	11.327	178	1119104	42.90	ng	99
73) Carbazole	11.486	167	991554	42.75	ng	100
74) Di-n-butylphthalate	11.804	149	1252099	46.93	ng	100
75) Fluoranthene	12.457	202	1160396	43.03	ng	99
77) Benzidine	12.586	184	552225	52.89	ng	99
78) Pyrene	12.692	202	1175089	45.53	ng	99
80) Butylbenzylphthalate	13.298	149	519359	50.74	ng	97
81) Benzo(a)anthracene	13.874	228	1007337	45.39	ng	99
82) 3,3'-Dichlorobenzidine	13.839	252	232198	28.21	ng	99
83) Chrysene	13.916	228	977227	44.85	ng	99
84) Bis(2-ethylhexyl)phtha...	13.857	149	740574	53.30	ng	99
85) Di-n-octyl phthalate	14.468	149	1208222	53.75	ng	100
87) Indeno(1,2,3-cd)pyrene	16.727	276	855975	45.68	ng	98
88) Benzo(b)fluoranthene	14.910	252	922756	47.30	ng	99
89) Benzo(k)fluoranthene	14.933	252	830498	44.81	ng	99
90) Benzo(a)pyrene	15.257	252	765458	45.42	ng	99
91) Dibenzo(a,h)anthracene	16.727	278	707269	45.83	ng	99
92) Benzo(g,h,i)perylene	17.139	276	708436	45.87	ng	98

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

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