

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050124\
 Data File : BF137792.D
 Acq On : 01 May 2024 15:21
 Operator : CG\JU
 Sample : P2306-02MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-4-20240429MS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/02/2024
 Supervised By :mohammad ahmed 05/03/2024

Quant Time: May 01 15:54:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:25:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.057	152	192854	20.000	ng	0.00
21) Naphthalene-d8	8.345	136	775329	20.000	ng	0.00
39) Acenaphthene-d10	10.110	164	440209	20.000	ng	0.00
64) Phenanthrene-d10	11.604	188	767868	20.000	ng	0.00
76) Chrysene-d12	14.257	240	434720	20.000	ng	# 0.00
86) Perylene-d12	15.827	264	441867	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	759154	65.907	ng	0.00
7) Phenol-d6	6.663	99	603393	41.299	ng	-0.01
23) Nitrobenzene-d5	7.622	82	1253138	93.902	ng	0.00
42) 2,4,6-Tribromophenol	10.904	330	624284	138.895	ng	0.00
45) 2-Fluorobiphenyl	9.422	172	2426931	89.443	ng	0.00
79) Terphenyl-d14	13.192	244	2681073	103.356	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.010	88	99984	19.276	ng	99
3) Pyridine	3.769	79	240355	19.146	ng	97
4) n-Nitrosodimethylamine	3.722	42	121766	21.540	ng	98
6) Aniline	6.716	93	436413	30.137	ng	99
8) 2-Chlorophenol	6.840	128	495773	39.495	ng	99
9) Benzaldehyde	6.610	77	361380	45.931	ng	96
10) Phenol	6.681	94	221945	14.825	ng	97
11) bis(2-Chloroethyl)ether	6.787	93	535809	45.709	ng	100
12) 1,3-Dichlorobenzene	6.998	146	517928	36.589	ng	99
13) 1,4-Dichlorobenzene	7.075	146	527671	37.150	ng	99
14) 1,2-Dichlorobenzene	7.228	146	522059	39.024	ng	98
15) Benzyl Alcohol	7.192	79	350694	34.710	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.322	45	720733	43.743	ng	99
17) 2-Methylphenol	7.298	107	334519	33.026	ng	98
18) Hexachloroethane	7.575	117	171598	35.737	ng	96
19) n-Nitroso-di-n-propyla...	7.463	70	413415	47.349	ng	98
20) 3+4-Methylphenols	7.445	107	396144	29.727	ng	97
22) Acetophenone	7.463	105	829393	47.564	ng	99
24) Nitrobenzene	7.639	77	621489	45.185	ng	99
25) Isophorone	7.875	82	1134267	47.141	ng	100
26) 2-Nitrophenol	7.957	139	321111	51.555	ng	99
27) 2,4-Dimethylphenol	7.981	122	400271	44.639	ng	97
28) bis(2-Chloroethoxy)met...	8.081	93	687066	47.161	ng	99
29) 2,4-Dichlorophenol	8.192	162	488133	45.038	ng	100
30) 1,2,4-Trichlorobenzene	8.281	180	480382	40.658	ng	99
31) Naphthalene	8.369	128	1657947	43.029	ng	100
32) Benzoic acid	8.045	122	100256	12.960	ng	99
33) 4-Chloroaniline	8.404	127	307361	21.512	ng	98
34) Hexachlorobutadiene	8.475	225	265729	36.632	ng	98
35) Caprolactam	8.775	113	32668m	9.513	ng	
36) 4-Chloro-3-methylphenol	8.875	107	474309	41.965	ng	97
37) 2-Methylnaphthalene	9.057	142	1131079	45.350	ng	99
38) 1-Methylnaphthalene	9.157	142	1064514	43.383	ng	98
40) 1,2,4,5-Tetrachloroben...	9.222	216	538568	46.193	ng	99
41) Hexachlorocyclopentadiene	9.210	237	542469	114.549	ng	99
43) 2,4,6-Trichlorophenol	9.328	196	383381	48.306	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.369	196	410362	46.827	ng	99
46) 1,1'-Biphenyl	9.522	154	1415484	46.854	ng	99
47) 2-Chloronaphthalene	9.551	162	1106172	45.352	ng	99
48) 2-Nitroaniline	9.645	65	337172	50.971	ng	92
49) Acenaphthylene	9.969	152	1792383	50.800	ng	99
50) Dimethylphthalate	9.822	163	1347701	48.496	ng	99
51) 2,6-Dinitrotoluene	9.886	165	310009	48.814	ng	88
52) Acenaphthene	10.145	154	1198684	51.254	ng	98
53) 3-Nitroaniline	10.051	138	175871	26.444	ng	95
54) 2,4-Dinitrophenol	10.163	184	329409	105.561	ng	98
55) Dibenzofuran	10.316	168	1613939	47.422	ng	98
56) 4-Nitrophenol	10.198	139	167221	30.337	ng	98
57) 2,4-Dinitrotoluene	10.292	165	420153	50.639	ng	88
58) Fluorene	10.657	166	1276828	47.140	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.428	232	349955	49.188	ng	99
60) Diethylphthalate	10.522	149	1252353	47.449	ng	99
61) 4-Chlorophenyl-phenyle...	10.645	204	624011	46.757	ng	98
62) 4-Nitroaniline	10.675	138	269578	39.962	ng	96
63) Azobenzene	10.810	77	1213666	47.577	ng	95
65) 4,6-Dinitro-2-methylph...	10.698	198	245294	55.656	ng	91
66) n-Nitrosodiphenylamine	10.769	169	1104983	48.074	ng	100
67) 4-Bromophenyl-phenylether	11.139	248	397898	48.419	ng	97
68) Hexachlorobenzene	11.210	284	428274	47.689	ng	# 93
69) Atrazine	11.292	200	364941	52.072	ng	98
70) Pentachlorophenol	11.398	266	555011	90.018	ng	99
71) Phenanthrene	11.627	178	1825482	48.774	ng	99
72) Anthracene	11.680	178	1857825	49.888	ng	100
73) Carbazole	11.833	167	1673346	46.625	ng	100
74) Di-n-butylphthalate	12.151	149	1895007	48.404	ng	100
75) Fluoranthene	12.821	202	1846395	45.131	ng	100
77) Benzidine	12.939	184	220234	21.905	ng	97
78) Pyrene	13.057	202	1857374	53.887	ng	100
80) Butylbenzylphthalate	13.663	149	657051	60.081	ng	93
81) Benzo(a)anthracene	14.245	228	1365380	49.107	ng	99
82) 3,3'-Dichlorobenzidine	14.198	252	257516	30.937	ng	99
83) Chrysene	14.280	228	1277130	49.056	ng	99
84) Bis(2-ethylhexyl)phtha...	14.210	149	874615	63.458	ng	100
85) Di-n-octyl phthalate	14.845	149	1280606	69.040	ng	99
87) Indeno(1,2,3-cd)pyrene	17.480	276	1389325	55.304	ng	99
88) Benzo(b)fluoranthene	15.357	252	1220124	44.381	ng	99
89) Benzo(k)fluoranthene	15.386	252	1213666	45.675	ng	100
90) Benzo(a)pyrene	15.762	252	1132663	50.501	ng	99
91) Dibenzo(a,h)anthracene	17.498	278	1100894	53.838	ng	98
92) Benzo(g,h,i)perylene	17.980	276	1095896	51.124	ng	98

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

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