

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052324\
 Data File : BF138029.D
 Acq On : 23 May 2024 16:00
 Operator : CG\JU
 Sample : P2556-02MSD
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-1MSD

Quant Time: May 23 16:25:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF051424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 15 04:12:16 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.033	152	144436	20.000	ng	-0.01
21) Naphthalene-d8	8.322	136	566520	20.000	ng	0.00
39) Acenaphthene-d10	10.086	164	314179	20.000	ng	0.00
64) Phenanthrene-d10	11.580	188	507545	20.000	ng	0.00
76) Chrysene-d12	14.233	240	260493	20.000	ng	0.00
86) Perylene-d12	15.792	264	296156	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	1056247	124.220	ng	0.02
7) Phenol-d6	6.663	99	1252852	119.611	ng	0.00
23) Nitrobenzene-d5	7.598	82	850716	88.373	ng	-0.01
42) 2,4,6-Tribromophenol	10.880	330	464221	146.541	ng	0.00
45) 2-Fluorobiphenyl	9.398	172	1835562	89.996	ng	0.00
79) Terphenyl-d14	13.162	244	1649744	93.338	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.169	88	136053	36.935	ng	# 86
3) Pyridine	3.845	79	313600	37.942	ng	99
4) n-Nitrosodimethylamine	3.793	42	165396	45.361	ng	98
6) Aniline	6.698	93	187992m	20.066	ng	
8) 2-Chlorophenol	6.822	128	437647	47.560	ng	99
9) Benzaldehyde	6.592	77	55697	9.828	ng	99
10) Phenol	6.675	94	457303	43.254	ng	96
11) bis(2-Chloroethyl)ether	6.769	93	379770	46.783	ng	98
12) 1,3-Dichlorobenzene	6.981	146	439712	41.300	ng	99
13) 1,4-Dichlorobenzene	7.051	146	450549	42.160	ng	98
14) 1,2-Dichlorobenzene	7.204	146	434335	43.183	ng	99
15) Benzyl Alcohol	7.175	79	367527	52.547	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.298	45	462487	46.349	ng	98
17) 2-Methylphenol	7.281	107	346581	47.228	ng	99
18) Hexachloroethane	7.545	117	154909	42.836	ng	96
19) n-Nitroso-di-n-propyla...	7.445	70	288572	49.037	ng	99
20) 3+4-Methylphenols	7.428	107	441396	45.912	ng	94
22) Acetophenone	7.445	105	588955	46.407	ng	# 99
24) Nitrobenzene	7.616	77	441898	45.250	ng	96
25) Isophorone	7.857	82	800920	47.938	ng	96
26) 2-Nitrophenol	7.933	139	239171	48.741	ng	99
27) 2,4-Dimethylphenol	7.957	122	326762	51.723	ng	98
28) bis(2-Chloroethoxy)met...	8.057	93	484240	47.822	ng	99
29) 2,4-Dichlorophenol	8.169	162	390527	49.158	ng	98
30) 1,2,4-Trichlorobenzene	8.257	180	393288	44.339	ng	99
31) Naphthalene	8.339	128	1268064	44.517	ng	99
32) Benzoic acid	8.098	122	315564	59.090	ng	95
33) 4-Chloroaniline	8.380	127	67862	7.267	ng	99
34) Hexachlorobutadiene	8.451	225	239291	42.621	ng	100
35) Caprolactam	8.775	113	114304m	49.663	ng	
36) 4-Chloro-3-methylphenol	8.869	107	397654	49.205	ng	99
37) 2-Methylnaphthalene	9.033	142	872621	47.658	ng	99
38) 1-Methylnaphthalene	9.133	142	801859	44.546	ng	100
40) 1,2,4,5-Tetrachloroben...	9.198	216	414457	48.523	ng	99
41) Hexachlorocyclopentadiene	9.186	237	359329	110.716	ng	98
43) 2,4,6-Trichlorophenol	9.310	196	289067	51.616	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.357	196	300784	49.086	ng	97
46) 1,1'-Biphenyl	9.498	154	1063137	48.245	ng	99
47) 2-Chloronaphthalene	9.527	162	827292	47.030	ng	98
48) 2-Nitroaniline	9.622	65	223057	49.518	ng	99
49) Acenaphthylene	9.951	152	1307344	51.688	ng	100
50) Dimethylphthalate	9.798	163	1005503	51.037	ng	99
51) 2,6-Dinitrotoluene	9.863	165	220530	48.771	ng	98
52) Acenaphthene	10.122	154	853966	51.083	ng	99
53) 3-Nitroaniline	10.033	138	92117	20.839	ng	98
54) 2,4-Dinitrophenol	10.139	184	194922	79.811	ng	# 86
55) Dibenzofuran	10.292	168	1183446	48.643	ng	100
56) 4-Nitrophenol	10.192	139	308377	97.128	ng	96
57) 2,4-Dinitrotoluene	10.269	165	297145	50.832	ng	# 88
58) Fluorene	10.639	166	927082	48.176	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.410	232	250008	51.935	ng	98
60) Diethylphthalate	10.504	149	938890	51.027	ng	99
61) 4-Chlorophenyl-phenyle...	10.621	204	463901	48.883	ng	99
62) 4-Nitroaniline	10.657	138	198145	47.259	ng	98
63) Azobenzene	10.786	77	832104	49.718	ng	99
65) 4,6-Dinitro-2-methylph...	10.680	198	141980	48.448	ng	97
66) n-Nitrosodiphenylamine	10.745	169	807656	51.890	ng	99
67) 4-Bromophenyl-phenylether	11.116	248	290228	52.160	ng	97
68) Hexachlorobenzene	11.186	284	298788	48.579	ng	99
69) Atrazine	11.274	200	257667	60.624	ng	99
70) Pentachlorophenol	11.380	266	355238	98.601	ng	99
71) Phenanthrene	11.604	178	1234626	49.363	ng	99
72) Anthracene	11.657	178	1256934	50.650	ng	99
73) Carbazole	11.810	167	1059052	46.718	ng	99
74) Di-n-butylphthalate	12.127	149	1340078	53.267	ng	99
75) Fluoranthene	12.798	202	1082565	42.642	ng	98
77) Benzidine	12.915	184	40253	11.824	ng	# 95
78) Pyrene	13.033	202	1073247	47.106	ng	99
80) Butylbenzylphthalate	13.633	149	409791	59.630	ng	99
81) Benzo(a)anthracene	14.221	228	824074	50.003	ng	100
82) 3,3'-Dichlorobenzidine	14.174	252	160512	35.144	ng	96
83) Chrysene	14.257	228	769706	48.835	ng	99
84) Bis(2-ethylhexyl)phtha...	14.186	149	513242	59.411	ng	99
85) Di-n-octyl phthalate	14.815	149	830809	71.189	ng	100
87) Indeno(1,2,3-cd)pyrene	17.421	276	726910	38.204	ng	99
88) Benzo(b)fluoranthene	15.321	252	841879	49.096	ng	99
89) Benzo(k)fluoranthene	15.356	252	807540	47.937	ng	100
90) Benzo(a)pyrene	15.727	252	763689	52.156	ng	99
91) Dibenzo(a,h)anthracene	17.445	278	596919	37.971	ng	98
92) Benzo(g,h,i)perylene	17.921	276	501440	30.364	ng	99

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

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