

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
 Data File : BF136271.D
 Acq On : 29 Nov 2023 02:57
 Operator : CG\JU
 Sample : 05553-01MS 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-01-112723MS

Quant Time: Nov 29 03:29:26 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 03:04:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.804	152	97573	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	344893	20.000	ng	0.00
39) Acenaphthene-d10	9.839	164	152601	20.000	ng	0.00
64) Phenanthrene-d10	11.328	188	317172	20.000	ng	0.00
76) Chrysene-d12	13.986	240	233200	20.000	ng	0.00
86) Perylene-d12	15.468	264	158895	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	132670	18.925	ng	0.00
7) Phenol-d6	6.434	99	146993	16.770	ng	-0.01
23) Nitrobenzene-d5	7.357	82	85887	13.159	ng	-0.02
42) 2,4,6-Tribromophenol	10.628	330	36831	19.741	ng	-0.01
45) 2-Fluorobiphenyl	9.157	172	157189	13.157	ng	-0.01
79) Terphenyl-d14	12.927	244	225893	11.173	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.563	88	23766	8.960	ng	98
3) Pyridine	3.334	79	49363	7.539	ng	96
4) n-Nitrosodimethylamine	3.269	42	28242	8.533	ng	# 94
6) Aniline	6.469	93	58463	7.057	ng	97
8) 2-Chlorophenol	6.587	128	66337	8.691	ng	97
9) Benzaldehyde	6.357	77	10256	2.099	ng	92
10) Phenol	6.445	94	80770	8.589	ng	91
11) bis(2-Chloroethyl)ether	6.540	93	64426	9.339	ng	97
12) 1,3-Dichlorobenzene	6.745	146	74218	8.568	ng	97
13) 1,4-Dichlorobenzene	6.822	146	74957	8.512	ng	95
14) 1,2-Dichlorobenzene	6.975	146	70766	8.488	ng	96
15) Benzyl Alcohol	6.940	79	50497	8.217	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.081	45	85686	8.942	ng	99
17) 2-Methylphenol	7.051	107	50074	8.344	ng	96
18) Hexachloroethane	7.316	117	22263	7.208	ng	# 84
19) n-Nitroso-di-n-propyla...	7.204	70	45710	8.719	ng	98
20) 3+4-Methylphenols	7.204	107	62506	8.014	ng	# 85
22) Acetophenone	7.210	105	95346	10.526	ng	# 86
24) Nitrobenzene	7.381	77	62391	9.433	ng	96
25) Isophorone	7.616	82	114035	9.690	ng	96
26) 2-Nitrophenol	7.698	139	25738	9.133	ng	99
27) 2,4-Dimethylphenol	7.734	122	44332	11.952	ng	94
28) bis(2-Chloroethoxy)met...	7.834	93	71586	10.056	ng	97
29) 2,4-Dichlorophenol	7.934	162	42914	8.578	ng	91
30) 1,2,4-Trichlorobenzene	8.022	180	55172	9.007	ng	98
31) Naphthalene	8.104	128	177933	9.179	ng	98
32) Benzoic acid	7.792	122	23755	7.027	ng	97
33) 4-Chloroaniline	8.151	127	45646	6.927	ng	98
34) Hexachlorobutadiene	8.222	225	33936	9.470	ng	97
35) Caprolactam	8.487	113	13111	9.374	ng	88
36) 4-Chloro-3-methylphenol	8.628	107	44648	8.228	ng	94
37) 2-Methylnaphthalene	8.792	142	104183	8.567	ng	98
38) 1-Methylnaphthalene	8.892	142	101370	8.545	ng	91
40) 1,2,4,5-Tetrachloroben...	8.957	216	50788	10.458	ng	97
41) Hexachlorocyclopentadiene	8.945	237	6304	3.373	ng	89
43) 2,4,6-Trichlorophenol	9.069	196	29398	9.326	ng	94

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112823\
 Data File : BF136271.D
 Acq On : 29 Nov 2023 02:57
 Operator : CG\JU
 Sample : 05553-01MS 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-01-112723MS

Quant Time: Nov 29 03:29:26 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 03:04:20 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.110	196	31757	9.232	ng	97
46) 1,1'-Biphenyl	9.257	154	132904	9.876	ng	95
47) 2-Chloronaphthalene	9.281	162	98170	9.394	ng	98
48) 2-Nitroaniline	9.375	65	25708	9.422	ng	97
49) Acenaphthylene	9.698	152	153502	10.293	ng	98
50) Dimethylphthalate	9.557	163	115450	10.010	ng	98
51) 2,6-Dinitrotoluene	9.622	165	21234	9.012	ng	92
52) Acenaphthene	9.869	154	90973	9.653	ng	91
53) 3-Nitroaniline	9.786	138	21335	9.125	ng	96
55) Dibenzofuran	10.045	168	134938	9.622	ng	96
56) 4-Nitrophenol	9.945	139	36444	20.524	ng	88
57) 2,4-Dinitrotoluene	10.022	165	26829	8.524	ng	93
58) Fluorene	10.386	166	110961	10.087	ng	94
59) 2,3,4,6-Tetrachlorophenol	10.163	232	28384	10.370	ng	97
60) Diethylphthalate	10.263	149	111717	9.683	ng	97
61) 4-Chlorophenyl-phenyle...	10.381	204	53957	10.100	ng	98
62) 4-Nitroaniline	10.398	138	24660	10.775	ng	94
63) Azobenzene	10.539	77	100256	9.373	ng	98
66) n-Nitrosodiphenylamine	10.498	169	94120	9.260	ng	98
67) 4-Bromophenyl-phenylether	10.875	248	31997	8.647	ng	93
68) Hexachlorobenzene	10.933	284	37054	8.652	ng	# 90
69) Atrazine	11.028	200	33838	11.223	ng	94
70) Pentachlorophenol	11.133	266	42537	17.572	ng	97
71) Phenanthrene	11.351	178	185081	10.382	ng	98
72) Anthracene	11.404	178	186294	10.421	ng	97
73) Carbazole	11.563	167	159180	10.469	ng	99
74) Di-n-butylphthalate	11.898	149	202946	10.668	ng	97
75) Fluoranthene	12.545	202	220282	12.172	ng	93
77) Benzidine	12.674	184	61796	16.120	ng	98
78) Pyrene	12.774	202	227445	8.906	ng	95
80) Butylbenzylphthalate	13.410	149	90502	11.386	ng	98
81) Benzo(a)anthracene	13.974	228	171199	9.884	ng	96
82) 3,3'-Dichlorobenzidine	13.939	252	44202	10.601	ng	98
83) Chrysene	14.010	228	158265	9.463	ng	94
84) Bis(2-ethylhexyl)phtha...	13.974	149	133472	13.992	ng	# 100
85) Di-n-octyl phthalate	14.592	149	204802	16.960	ng	# 95
87) Indeno(1,2,3-cd)pyrene	16.968	276	94870	10.221	ng	96
88) Benzo(b)fluoranthene	15.027	252	112366	10.080	ng	# 94
89) Benzo(k)fluoranthene	15.057	252	101409	9.578	ng	# 94
90) Benzo(a)pyrene	15.404	252	87202	9.569	ng	# 91
91) Dibenzo(a,h)anthracene	16.992	278	76242	10.259	ng	# 90
92) Benzo(g,h,i)perylene	17.421	276	78297	9.762	ng	95

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

