

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081422\
 Data File : BF129761.D
 Acq On : 14 Aug 2022 14:06
 Operator : CG\JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC005

Quant Time: Aug 15 01:30:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 15 01:28:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	156313	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	627416	20.000	ng	0.00
39) Acenaphthene-d10	9.863	164	340110	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	611475	20.000	ng	0.00
76) Chrysene-d12	13.998	240	486221	20.000	ng	0.00
86) Perylene-d12	15.468	264	393168	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	104281	10.868	ng	0.00
7) Phenol-d6	6.481	99	132778	11.009	ng	0.00
23) Nitrobenzene-d5	7.387	82	120521	10.486	ng	-0.01
42) 2,4,6-Tribromophenol	10.657	330	35349	10.810	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	247166	11.242	ng	-0.01
79) Terphenyl-d14	12.933	244	283890	11.015	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.581	88	20728	4.788	ng	96
3) Pyridine	3.440	79	46719	3.886	ng	99
4) n-Nitrosodimethylamine	3.310	42	23896	4.526	ng	# 98
6) Aniline	6.487	93	74208	4.949	ng	# 73
8) 2-Chlorophenol	6.616	128	57806	5.514	ng	96
10) Phenol	6.492	94	70510	5.490	ng	95
11) bis(2-Chloroethyl)ether	6.557	93	51657	5.071	ng	97
12) 1,3-Dichlorobenzene	6.763	146	61877	5.503	ng	95
13) 1,4-Dichlorobenzene	6.839	146	63427	5.547	ng	97
14) 1,2-Dichlorobenzene	6.992	146	59729	5.683	ng	94
15) Benzyl Alcohol	6.969	79	47978	5.485	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.092	45	72528	4.737	ng	85
17) 2-Methylphenol	7.092	107	46223	5.315	ng	96
18) Hexachloroethane	7.328	117	23485	5.454	ng	95
19) n-Nitroso-di-n-propyla...	7.228	70	39893	5.464	ng	98
20) 3+4-Methylphenols	7.245	107	62743	5.674	ng	95
22) Acetophenone	7.234	105	81629	5.588	ng	# 94
24) Nitrobenzene	7.404	77	62216	5.257	ng	97
25) Isophorone	7.639	82	103308	5.014	ng	97
26) 2-Nitrophenol	7.728	139	27892	4.777	ng	95
27) 2,4-Dimethylphenol	7.769	122	43057	4.916	ng	98
28) bis(2-Chloroethoxy)met...	7.851	93	61073	5.110	ng	99
29) 2,4-Dichlorophenol	7.981	162	46067	5.096	ng	96
30) 1,2,4-Trichlorobenzene	8.045	180	50698	5.418	ng	99
31) Naphthalene	8.128	128	176096	5.524	ng	99
33) 4-Chloroaniline	8.181	127	67317	5.144	ng	97
34) Hexachlorobutadiene	8.234	225	30242	5.686	ng	99
35) Caprolactam	8.528	113	14393	4.897	ng	99
36) 4-Chloro-3-methylphenol	8.675	107	50497	5.267	ng	98
37) 2-Methylnaphthalene	8.816	142	113014	5.456	ng	95
38) 1-Methylnaphthalene	8.916	142	112433	5.622	ng	99
40) 1,2,4,5-Tetrachloroben...	8.986	216	49212	5.511	ng	98
43) 2,4,6-Trichlorophenol	9.104	196	30382	4.684	ng	98
44) 2,4,5-Trichlorophenol	9.157	196	32284	4.577	ng	98
46) 1,1'-Biphenyl	9.281	154	139381	5.615	ng	98
47) 2-Chloronaphthalene	9.310	162	108289	5.397	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081422\
 Data File : BF129761.D
 Acq On : 14 Aug 2022 14:06
 Operator : CG\JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC005

Quant Time: Aug 15 01:30:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 15 01:28:08 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
48) 2-Nitroaniline	9.410	65	32138	4.817	ng	98
49) Acenaphthylene	9.722	152	166736	5.469	ng	98
50) Dimethylphthalate	9.575	163	129649	5.522	ng	99
51) 2,6-Dinitrotoluene	9.645	165	26694	5.080	ng	95
52) Acenaphthene	9.892	154	100468	5.045	ng	99
53) 3-Nitroaniline	9.822	138	30223	4.870	ng	98
55) Dibenzofuran	10.063	168	155652	5.670	ng	98
56) 4-Nitrophenol	10.022	139	9325	2.037	ng	90
57) 2,4-Dinitrotoluene	10.057	165	36397	5.302	ng #	92
58) Fluorene	10.410	166	130094	5.923	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.198	232	22463	4.121	ng #	78
60) Diethylphthalate	10.275	149	131194	5.780	ng	98
61) 4-Chlorophenyl-phenyle...	10.398	204	56464	5.831	ng	96
62) 4-Nitroaniline	10.433	138	30417	4.941	ng	99
63) Azobenzene	10.557	77	122054	5.272	ng	99
65) 4,6-Dinitro-2-methylph...	10.475	198	13422	3.478	ng	97
66) n-Nitrosodiphenylamine	10.516	169	106134	5.212	ng	98
67) 4-Bromophenyl-phenylether	10.886	248	36223	5.410	ng	95
68) Hexachlorobenzene	10.957	284	39508	5.446	ng	99
69) Atrazine	11.045	200	33926	5.485	ng	99
71) Phenanthrene	11.374	178	182823	5.477	ng	99
72) Anthracene	11.427	178	182496	5.564	ng	98
73) Carbazole	11.586	167	168808	5.467	ng	99
74) Di-n-butylphthalate	11.904	149	213567	5.808	ng	98
75) Fluoranthene	12.569	202	194658	5.756	ng	99
78) Pyrene	12.798	202	197130	4.848	ng	99
80) Butylbenzylphthalate	13.404	149	86879	4.926	ng	97
81) Benzo(a)anthracene	13.986	228	173994	5.239	ng	98
82) 3,3'-Dichlorobenzidine	13.951	252	58480	5.333	ng	98
83) Chrysene	14.027	228	165261	5.279	ng	97
84) Bis(2-ethylhexyl)phtha...	13.957	149	107894	4.647	ng	97
85) Di-n-octyl phthalate	14.568	149	147421	4.412	ng	99
87) Indeno(1,2,3-cd)pyrene	16.945	276	109859	4.149	ng	98
88) Benzo(b)fluoranthene	15.039	252	141927	5.442	ng	98
89) Benzo(k)fluoranthene	15.062	252	136887	5.356	ng	99
90) Benzo(a)pyrene	15.404	252	109679	5.181	ng	98
91) Dibenzo(a,h)anthracene	16.951	278	89252	4.142	ng	97
92) Benzo(g,h,i)perylene	17.392	276	84140	3.821	ng #	89

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

Instrument :

BNA F

ClientSampleId :

SSTDICC005

Quant Time: Aug 15 01:30:59 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Aug 15 01:28:08 2022

Response via : Initial Calibration

