

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010424\
 Data File : BF136962.D
 Acq On : 05 Jan 2024 02:31
 Operator : CG\JU
 Sample : P1004-01MSD 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-01022024MSD

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 01/05/2024
 Supervised By :mohammad ahmed 01/05/2024

Quant Time: Jan 05 03:55:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	46666	20.000	ng	-0.01
21) Naphthalene-d8	8.057	136	171726	20.000	ng	-0.01
39) Acenaphthene-d10	9.816	164	82066	20.000	ng	-0.01
64) Phenanthrene-d10	11.310	188	146476	20.000	ng	0.00
76) Chrysene-d12	13.969	240	112186	20.000	ng	0.00
86) Perylene-d12	15.457	264	85036	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	165419	55.308	ng	0.00
7) Phenol-d6	6.434	99	199176	51.394	ng	0.00
23) Nitrobenzene-d5	7.345	82	117850	35.117	ng	-0.01
42) 2,4,6-Tribromophenol	10.610	330	48860	50.553	ng	-0.01
45) 2-Fluorobiphenyl	9.134	172	249987	40.971	ng	-0.02
79) Terphenyl-d14	12.904	244	251486	31.327	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.499	88	23952	19.220	ng	96
3) Pyridine	3.275	79	53628	17.638	ng	88
4) n-Nitrosodimethylamine	3.222	42	30425	18.738	ng	# 93
6) Aniline	6.451	93	69909	18.043	ng	# 63
8) 2-Chlorophenol	6.575	128	69976	21.380	ng	94
9) Benzaldehyde	6.340	77	14535	6.594	ng	96
10) Phenol	6.445	94	87794	21.221	ng	91
11) bis(2-Chloroethyl)ether	6.522	93	63882	20.305	ng	93
12) 1,3-Dichlorobenzene	6.722	146	71396	19.996	ng	97
13) 1,4-Dichlorobenzene	6.798	146	76118	20.864	ng	99
14) 1,2-Dichlorobenzene	6.951	146	68737	20.248	ng	97
15) Benzyl Alcohol	6.928	79	53119	20.317	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.057	45	100544	19.547	ng	76
17) 2-Methylphenol	7.051	107	53178	19.612	ng	# 90
18) Hexachloroethane	7.287	117	22606	17.062	ng	92
19) n-Nitroso-di-n-propyla...	7.192	70	47087	20.043	ng	95
20) 3+4-Methylphenols	7.204	107	71481	21.102	ng	# 52
22) Acetophenone	7.192	105	99623	23.145	ng	# 85
24) Nitrobenzene	7.363	77	69158	20.572	ng	99
25) Isophorone	7.598	82	125106	20.488	ng	99
26) 2-Nitrophenol	7.681	139	31325	19.476	ng	100
27) 2,4-Dimethylphenol	7.722	122	52764	26.946	ng	98
28) bis(2-Chloroethoxy)met...	7.810	93	80286	21.630	ng	98
29) 2,4-Dichlorophenol	7.934	162	52869	20.972	ng	94
30) 1,2,4-Trichlorobenzene	8.004	180	59796	21.289	ng	99
31) Naphthalene	8.081	128	199885	21.187	ng	100
32) Benzoic acid	7.828	122	31683	16.722	ng	94
33) 4-Chloroaniline	8.139	127	63840	18.328	ng	94
34) Hexachlorobutadiene	8.192	225	34908	20.456	ng	98
35) Caprolactam	8.487	113	17334	20.843	ng	96
36) 4-Chloro-3-methylphenol	8.628	107	58021	20.444	ng	95
37) 2-Methylnaphthalene	8.775	142	124147	20.447	ng	98
38) 1-Methylnaphthalene	8.875	142	120821	20.283	ng	95
40) 1,2,4,5-Tetrachloroben...	8.939	216	61114	24.046	ng	98
41) Hexachlorocyclopentadiene	8.922	237	467	5.439	ng	78
43) 2,4,6-Trichlorophenol	9.063	196	36026	22.103	ng	97

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44) 2,4,5-Trichlorophenol	9.110	196	39567	21.784	ng	97
46) 1,1'-Biphenyl	9.239	154	163619	23.759	ng	99
47) 2-Chloronaphthalene	9.263	162	115426	22.048	ng	98
48) 2-Nitroaniline	9.369	65	36278	22.129	ng	87
49) Acenaphthylene	9.681	152	186290	23.280	ng	99
50) Dimethylphthalate	9.539	163	132691	22.136	ng	99
51) 2,6-Dinitrotoluene	9.610	165	27013	19.856	ng	92
52) Acenaphthene	9.851	154	108411	21.856	ng	98
53) 3-Nitroaniline	9.781	138	31059	21.552	ng	98
54) 2,4-Dinitrophenol	9.904	184	1713	5.901	ng	# 89
55) Dibenzofuran	10.022	168	166535	22.148	ng	99
56) 4-Nitrophenol	9.969	139	31489	32.368	ng	95
57) 2,4-Dinitrotoluene	10.016	165	32742	18.539	ng	# 98
58) Fluorene	10.369	166	132456	22.201	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.151	232	27773	19.864	ng	99
60) Diethylphthalate	10.239	149	135556	21.795	ng	99
61) 4-Chlorophenyl-phenyle...	10.357	204	62200	21.449	ng	98
62) 4-Nitroaniline	10.392	138	29706m	21.383	ng	
63) Azobenzene	10.522	77	116629	20.077	ng	98
65) 4,6-Dinitro-2-methylph...	10.433	198	2450	2.919	ng	74
66) n-Nitrosodiphenylamine	10.481	169	114466	24.057	ng	96
67) 4-Bromophenyl-phenylether	10.851	248	36253	22.286	ng	98
68) Hexachlorobenzene	10.916	284	38816	21.412	ng	94
69) Atrazine	11.010	200	35720	29.351	ng	99
70) Pentachlorophenol	11.122	266	28902	34.204	ng	98
71) Phenanthrene	11.333	178	185989	24.746	ng	99
72) Anthracene	11.386	178	182025	23.915	ng	98
73) Carbazole	11.551	167	166944	23.263	ng	99
74) Di-n-butylphthalate	11.875	149	212134	24.225	ng	98
75) Fluoranthene	12.533	202	230854	28.716	ng	99
77) Benzidine	12.663	184	70753	21.389	ng	99
78) Pyrene	12.763	202	230475	20.928	ng	99
80) Butylbenzylphthalate	13.386	149	85789	21.408	ng	97
81) Benzo(a)anthracene	13.963	228	198982	25.543	ng	98
82) 3,3'-Dichlorobenzidine	13.927	252	57577	26.026	ng	94
83) Chrysene	13.998	228	185615	24.366	ng	96
84) Bis(2-ethylhexyl)phtha...	13.945	149	130036	25.791	ng	98
85) Di-n-octyl phthalate	14.563	149	217304	27.863	ng	98
87) Indeno(1,2,3-cd)pyrene	16.980	276	118401	19.284	ng	98
88) Benzo(b)fluoranthene	15.021	252	150726	26.538	ng	98
89) Benzo(k)fluoranthene	15.051	252	136169	25.373	ng	98
90) Benzo(a)pyrene	15.398	252	115206	24.030	ng	98
91) Dibenzo(a,h)anthracene	16.986	278	98514	19.558	ng	98
92) Benzo(g,h,i)perylene	17.433	276	94597	18.336	ng	100

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

