

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072423\
 Data File : BF134430.D
 Acq On : 24 Jul 2023 15:28
 Operator : CG\JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC020

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/25/2023
 Supervised By :mohammad ahmed 07/25/2023

Quant Time: Jul 25 01:35:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 01:30:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.833	152	48428	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	179271	20.000	ng	0.00
39) Acenaphthene-d10	9.874	164	88329	20.000	ng	0.00
64) Phenanthrene-d10	11.368	188	178575	20.000	ng	0.00
76) Chrysene-d12	14.033	240	148098	20.000	ng	0.00
86) Perylene-d12	15.533	264	130938	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	126947	43.310	ng	0.00
7) Phenol-d6	6.475	99	156440	43.562	ng	0.00
23) Nitrobenzene-d5	7.398	82	145831	41.157	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	53061	42.280	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	283403	42.121	ng	0.00
79) Terphenyl-d14	12.962	244	345437	40.717	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.593	88	25930	20.856	ng	96
3) Pyridine	3.387	79	58351m	20.944	ng	
4) n-Nitrosodimethylamine	3.334	42	35750	21.676	ng	98
6) Aniline	6.498	93	91505	21.619	ng	94
8) 2-Chlorophenol	6.622	128	63751	21.469	ng	97
9) Benzaldehyde	6.392	77	43343	22.537	ng	97
10) Phenol	6.486	94	80543	21.556	ng	94
11) bis(2-Chloroethyl)ether	6.575	93	52852	20.262	ng	94
12) 1,3-Dichlorobenzene	6.775	146	69312	21.706	ng	95
13) 1,4-Dichlorobenzene	6.851	146	70225	21.370	ng	99
14) 1,2-Dichlorobenzene	7.004	146	66777	21.633	ng	97
15) Benzyl Alcohol	6.981	79	60133	21.901	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.110	45	80707	21.010	ng	99
17) 2-Methylphenol	7.092	107	56119	21.191	ng	99
18) Hexachloroethane	7.339	117	27177	21.527	ng	98
19) n-Nitroso-di-n-propyla...	7.239	70	47451	21.225	ng	98
20) 3+4-Methylphenols	7.245	107	73029	21.538	ng	92
22) Acetophenone	7.245	105	92799	20.814	ng	# 91
24) Nitrobenzene	7.416	77	72190	20.486	ng	98
25) Isophorone	7.651	82	120857	20.466	ng	97
26) 2-Nitrophenol	7.733	139	34025	20.993	ng	95
27) 2,4-Dimethylphenol	7.769	122	53380	20.870	ng	95
28) bis(2-Chloroethoxy)met...	7.863	93	68024	20.695	ng	98
29) 2,4-Dichlorophenol	7.975	162	52169	20.404	ng	97
30) 1,2,4-Trichlorobenzene	8.057	180	55919	20.275	ng	95
31) Naphthalene	8.133	128	184294	20.656	ng	99
32) Benzoic acid	7.869	122	35403m	20.961	ng	
33) 4-Chloroaniline	8.186	127	77184	20.929	ng	95
34) Hexachlorobutadiene	8.245	225	39060	21.141	ng	98
35) Caprolactam	8.545	113	15493	19.855	ng	96
36) 4-Chloro-3-methylphenol	8.669	107	59295	20.099	ng	98
37) 2-Methylnaphthalene	8.827	142	128273	20.692	ng	98
38) 1-Methylnaphthalene	8.927	142	120732	20.993	ng	93
40) 1,2,4,5-Tetrachloroben...	8.992	216	57745	20.629	ng	100
41) Hexachlorocyclopentadiene	8.975	237	9782	19.775	ng	90
43) 2,4,6-Trichlorophenol	9.110	196	37069	20.643	ng	94

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44) 2,4,5-Trichlorophenol	9.157	196	44452	21.247	ng	98
46) 1,1'-Biphenyl	9.292	154	153247	21.050	ng	99
47) 2-Chloronaphthalene	9.316	162	111139	20.708	ng	97
48) 2-Nitroaniline	9.416	65	38955	20.154	ng	93
49) Acenaphthylene	9.733	152	190027	20.958	ng	98
50) Dimethylphthalate	9.592	163	136120	20.789	ng	99
51) 2,6-Dinitrotoluene	9.657	165	29294	20.389	ng	99
52) Acenaphthene	9.904	154	111676	20.850	ng	99
53) 3-Nitroaniline	9.833	138	35471	21.533	ng	99
54) 2,4-Dinitrophenol	9.951	184	10760	19.672	ng	96
55) Dibenzofuran	10.080	168	176383	21.396	ng	96
56) 4-Nitrophenol	10.010	139	16660	18.965	ng	87
57) 2,4-Dinitrotoluene	10.069	165	39793	20.621	ng	# 93
58) Fluorene	10.421	166	138096	20.765	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.198	232	31402m	19.462	ng	
60) Diethylphthalate	10.292	149	139654	20.600	ng	98
61) 4-Chlorophenyl-phenyle...	10.416	204	67610	21.282	ng	97
62) 4-Nitroaniline	10.445	138	35220	20.847	ng	91
63) Azobenzene	10.574	77	127572	20.444	ng	97
65) 4,6-Dinitro-2-methylph...	10.480	198	20249	19.768	ng	98
66) n-Nitrosodiphenylamine	10.533	169	118238	20.811	ng	97
67) 4-Bromophenyl-phenylether	10.904	248	40966	20.938	ng	# 90
68) Hexachlorobenzene	10.974	284	46051	20.902	ng	93
69) Atrazine	11.063	200	33413	21.094	ng	92
70) Pentachlorophenol	11.174	266	17137	18.534	ng	93
71) Phenanthrene	11.392	178	196683	21.080	ng	99
72) Anthracene	11.445	178	199792	20.747	ng	99
73) Carbazole	11.604	167	188724	20.812	ng	100
74) Di-n-butylphthalate	11.927	149	215165	20.670	ng	99
75) Fluoranthene	12.586	202	224371	20.634	ng	99
77) Benzidine	12.715	184	64456	20.383	ng	99
78) Pyrene	12.821	202	237451	20.165	ng	98
80) Butylbenzylphthalate	13.439	149	95741	20.297	ng	97
81) Benzo(a)anthracene	14.021	228	214727	20.782	ng	97
82) 3,3'-Dichlorobenzidine	13.980	252	71739	21.921	ng	96
83) Chrysene	14.057	228	211457	21.396	ng	98
84) Bis(2-ethylhexyl)phtha...	14.004	149	118543	20.686	ng	98
85) Di-n-octyl phthalate	14.621	149	160412	20.143	ng	100
87) Indeno(1,2,3-cd)pyrene	17.080	276	160054	18.710	ng	99
88) Benzo(b)fluoranthene	15.092	252	171812	20.974	ng	99
89) Benzo(k)fluoranthene	15.121	252	173963m	21.589	ng	
90) Benzo(a)pyrene	15.468	252	156739	20.701	ng	99
91) Dibenzo(a,h)anthracene	17.092	278	129831	18.793	ng	97
92) Benzo(g,h,i)perylene	17.550	276	136765	19.281	ng	99

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

