

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
 Data File : BF130283.D
 Acq On : 16 Sep 2022 10:12
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/19/2022
 Supervised By : Jagrut Upadhyay 09/19/2022

Quant Time: Sep 17 02:28:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 14 14:55:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.681	152	103515	20.000	ng	0.00
21) Naphthalene-d8	7.963	136	393538	20.000	ng	# 0.00
39) Acenaphthene-d10	9.710	164	208246	20.000	ng	0.00
64) Phenanthrene-d10	11.192	188	344735	20.000	ng	0.00
76) Chrysene-d12	13.821	240	188114	20.000	ng	0.00
86) Perylene-d12	15.209	264	155717	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.275	112	464040	77.753	ng	0.00
7) Phenol-d6	6.316	99	575169	77.023	ng	0.00
23) Nitrobenzene-d5	7.251	82	586020	76.076	ng	0.00
42) 2,4,6-Tribromophenol	10.498	330	169886	85.139	ng	0.00
45) 2-Fluorobiphenyl	9.039	172	1087107	76.018	ng	0.00
79) Terphenyl-d14	12.774	244	981898	86.464	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.299	88	96470	35.196	ng	97
3) Pyridine	2.987	79	265028	37.067	ng	97
4) n-Nitrosodimethylamine	2.951	42	131945	33.930	ng	95
6) Aniline	6.345	93	351717	38.226	ng	98
8) 2-Chlorophenol	6.463	128	266334	39.176	ng	93
9) Benzaldehyde	6.228	77	185911	37.167	ng	94
10) Phenol	6.328	94	330630	37.781	ng	99
11) bis(2-Chloroethyl)ether	6.422	93	240128	38.043	ng	96
12) 1,3-Dichlorobenzene	6.616	146	294559	39.376	ng	99
13) 1,4-Dichlorobenzene	6.698	146	298627	39.147	ng	98
14) 1,2-Dichlorobenzene	6.851	146	279714	39.252	ng	97
15) Benzyl Alcohol	6.828	79	219063	37.000	ng	96
16) 2,2'-oxybis(1-Chloropr...	6.963	45	324009	35.194	ng	97
17) 2-Methylphenol	6.939	107	212787	38.503	ng	99
18) Hexachloroethane	7.186	117	115183	38.305	ng	97
19) n-Nitroso-di-n-propyla...	7.104	70	189142	37.536	ng	94
20) 3+4-Methylphenols	7.092	107	276284	38.484	ng	88
22) Acetophenone	7.092	105	365039	37.940	ng	# 99
24) Nitrobenzene	7.269	77	296214	37.871	ng	95
25) Isophorone	7.504	82	476140	38.350	ng	99
26) 2-Nitrophenol	7.580	139	136561	42.591	ng	91
27) 2,4-Dimethylphenol	7.622	122	201417	40.798	ng	94
28) bis(2-Chloroethoxy)met...	7.722	93	271213	38.795	ng	99
29) 2,4-Dichlorophenol	7.822	162	221432	40.929	ng	96
30) 1,2,4-Trichlorobenzene	7.904	180	234850	40.151	ng	96
31) Naphthalene	7.986	128	788543	38.893	ng	99
32) Benzoic acid	7.745	122	158774m	41.587	ng	
33) 4-Chloroaniline	8.033	127	306613	39.763	ng	99
34) Hexachlorobutadiene	8.098	225	153254	40.997	ng	99
35) Caprolactam	8.416	113	65894m	42.486	ng	
36) 4-Chloro-3-methylphenol	8.516	107	240001	40.023	ng	99
37) 2-Methylnaphthalene	8.675	142	512988	39.684	ng	99
38) 1-Methylnaphthalene	8.775	142	500050	40.268	ng	99
40) 1,2,4,5-Tetrachloroben...	8.839	216	221158	38.915	ng	99
41) Hexachlorocyclopentadiene	8.822	237	118851	39.061	ng	99
43) 2,4,6-Trichlorophenol	8.951	196	156980	39.578	ng	93

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44) 2,4,5-Trichlorophenol	8.986	196	172433	40.287	ng	94
46) 1,1'-Biphenyl	9.139	154	592757	38.113	ng	98
47) 2-Chloronaphthalene	9.163	162	484463	38.507	ng	97
48) 2-Nitroaniline	9.263	65	157622	37.563	ng	87
49) Acenaphthylene	9.574	152	723720	38.366	ng	99
50) Dimethylphthalate	9.445	163	561459	39.031	ng	99
51) 2,6-Dinitrotoluene	9.504	165	122872	40.843	ng	90
52) Acenaphthene	9.745	154	477951m	38.563	ng	
53) 3-Nitroaniline	9.674	138	133918	39.949	ng	91
54) 2,4-Dinitrophenol	9.774	184	61728	41.122	ng	# 87
55) Dibenzofuran	9.916	168	666933	38.687	ng	99
56) 4-Nitrophenol	9.827	139	97350	39.872	ng	93
57) 2,4-Dinitrotoluene	9.904	165	161933	42.118	ng	87
58) Fluorene	10.257	166	550342	39.036	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.033	232	137643	41.086	ng	96
60) Diethylphthalate	10.139	149	563484	39.063	ng	100
61) 4-Chlorophenyl-phenyle...	10.257	204	244962	39.608	ng	93
62) 4-Nitroaniline	10.286	138	133993	39.790	ng	92
63) Azobenzene	10.416	77	541032	36.424	ng	94
65) 4,6-Dinitro-2-methylph...	10.310	198	84546	44.057	ng	84
66) n-Nitrosodiphenylamine	10.374	169	451347	39.180	ng	99
67) 4-Bromophenyl-phenylether	10.739	248	156272	41.092	ng	94
68) Hexachlorobenzene	10.798	284	178182	41.332	ng	94
69) Atrazine	10.904	200	132687	39.413	ng	97
70) Pentachlorophenol	10.992	266	98278	42.478	ng	98
71) Phenanthrene	11.216	178	757130	39.118	ng	99
72) Anthracene	11.269	178	737661	39.494	ng	100
73) Carbazole	11.427	167	642597	38.122	ng	99
74) Di-n-butylphthalate	11.763	149	787755	38.754	ng	99
75) Fluoranthene	12.398	202	710787	38.005	ng	98
77) Benzidine	12.527	184	212242	38.867	ng	99
78) Pyrene	12.627	202	704207	42.793	ng	100
80) Butylbenzylphthalate	13.251	149	252274	41.381	ng	99
81) Benzo(a)anthracene	13.810	228	488707	39.052	ng	99
82) 3,3'-Dichlorobenzidine	13.780	252	137008	40.219	ng	# 98
83) Chrysene	13.845	228	467498	39.344	ng	99
84) Bis(2-ethylhexyl)phtha...	13.810	149	281649	40.333	ng	99
85) Di-n-octyl phthalate	14.421	149	417090	39.051	ng	98
87) Indeno(1,2,3-cd)pyrene	16.545	276	532042	48.181	ng	98
88) Benzo(b)fluoranthene	14.821	252	391063	38.477	ng	99
89) Benzo(k)fluoranthene	14.845	252	383950	38.403	ng	100
90) Benzo(a)pyrene	15.151	252	327443	40.666	ng	99
91) Dibenzo(a,h)anthracene	16.568	278	436884	48.228	ng	97
92) Benzo(g,h,i)perylene	16.950	276	436729	47.859	ng	100

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

