

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122324\
 Data File : BF140961.D
 Acq On : 23 Dec 2024 10:18
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/24/2024
 Supervised By :mohammad ahmed 12/24/2024

Quant Time: Dec 23 13:23:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 16:59:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	61135	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	219500	20.000	ng	0.00
39) Acenaphthene-d10	9.886	164	122112	20.000	ng	0.00
64) Phenanthrene-d10	11.374	188	222339	20.000	ng	0.00
76) Chrysene-d12	14.033	240	134875	20.000	ng	0.00
86) Perylene-d12	15.533	264	144038	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	310519	83.614	ng	0.00
7) Phenol-d6	6.504	99	393458	79.989	ng	0.00
23) Nitrobenzene-d5	7.422	82	360941	82.270	ng	0.00
42) 2,4,6-Tribromophenol	10.680	330	112671	78.014	ng	0.00
45) 2-Fluorobiphenyl	9.204	172	755529	85.064	ng	0.00
79) Terphenyl-d14	12.957	244	735775	86.022	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.593	88	59611	37.265	ng	97
3) Pyridine	3.375	79	138053	37.266	ng	98
4) n-Nitrosodimethylamine	3.334	42	68250	36.546	ng	96
6) Aniline	6.522	93	162833	38.717	ng	95
8) 2-Chlorophenol	6.645	128	164578	40.638	ng	99
9) Benzaldehyde	6.410	77	96199	35.296	ng	98
10) Phenol	6.516	94	202012	39.625	ng	96
11) bis(2-Chloroethyl)ether	6.587	93	154671	40.682	ng	99
12) 1,3-Dichlorobenzene	6.792	146	191724	41.546	ng	97
13) 1,4-Dichlorobenzene	6.869	146	191959	40.980	ng	100
14) 1,2-Dichlorobenzene	7.022	146	184256	41.690	ng	99
15) Benzyl Alcohol	7.004	79	140053	39.822	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.122	45	176332	39.371	ng	85
17) 2-Methylphenol	7.122	107	129035	39.908	ng	98
18) Hexachloroethane	7.357	117	68878	39.498	ng	97
19) n-Nitroso-di-n-propyla...	7.263	70	113462	38.454	ng	99
20) 3+4-Methylphenols	7.275	107	173324	40.426	ng	# 85
22) Acetophenone	7.263	105	233626	42.548	ng	95
24) Nitrobenzene	7.439	77	183431	40.811	ng	96
25) Isophorone	7.675	82	301414	41.003	ng	99
26) 2-Nitrophenol	7.751	139	88301	42.656	ng	100
27) 2,4-Dimethylphenol	7.792	122	100833	41.709	ng	97
28) bis(2-Chloroethoxy)met...	7.875	93	191575	41.726	ng	98
29) 2,4-Dichlorophenol	8.004	162	142754	42.771	ng	98
30) 1,2,4-Trichlorobenzene	8.075	180	165052	43.138	ng	99
31) Naphthalene	8.151	128	504908	41.875	ng	100
32) Benzoic acid	7.945	122	88707	38.690	ng	97
33) 4-Chloroaniline	8.210	127	162948	40.690	ng	98
34) Hexachlorobutadiene	8.263	225	110624	42.974	ng	99
35) Caprolactam	8.592	113	38139	38.383	ng	94
36) 4-Chloro-3-methylphenol	8.698	107	149244	39.729	ng	99
37) 2-Methylnaphthalene	8.839	142	327151	41.017	ng	99
38) 1-Methylnaphthalene	8.939	142	323764	41.045	ng	100
40) 1,2,4,5-Tetrachloroben...	9.010	216	167089	44.518	ng	98
41) Hexachlorocyclopentadiene	8.986	237	22091	28.993	ng	95
43) 2,4,6-Trichlorophenol	9.133	196	94594	41.160	ng	99

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44) 2,4,5-Trichlorophenol	9.186	196	99644	39.447	ng	98
46) 1,1'-Biphenyl	9.304	154	416413	42.884	ng	99
47) 2-Chloronaphthalene	9.333	162	318010	42.810	ng	99
48) 2-Nitroaniline	9.439	65	88554	39.472	ng	94
49) Acenaphthylene	9.745	152	465833	42.672	ng	99
50) Dimethylphthalate	9.604	163	350137	40.670	ng	100
51) 2,6-Dinitrotoluene	9.681	165	80680	40.993	ng	91
52) Acenaphthene	9.916	154	297912	42.722	ng	99
53) 3-Nitroaniline	9.851	138	77647	39.755	ng	99
54) 2,4-Dinitrophenol	9.980	184	31040	36.222	ng	96
55) Dibenzofuran	10.092	168	447317	41.397	ng	98
56) 4-Nitrophenol	10.057	139	14441m	17.424	ng	
57) 2,4-Dinitrotoluene	10.086	165	105506	40.743	ng	98
58) Fluorene	10.433	166	361115	40.509	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.222	232	81536	39.612	ng	95
60) Diethylphthalate	10.298	149	358822	40.193	ng	99
61) 4-Chlorophenyl-phenyle...	10.422	204	186101	41.560	ng	97
62) 4-Nitroaniline	10.469	138	77257m	37.852	ng	
63) Azobenzene	10.580	77	327136	38.978	ng	99
65) 4,6-Dinitro-2-methylph...	10.504	198	52517	41.684	ng	96
66) n-Nitrosodiphenylamine	10.545	169	303009	44.326	ng	99
67) 4-Bromophenyl-phenylether	10.910	248	109503	44.024	ng	99
68) Hexachlorobenzene	10.986	284	124414	44.116	ng	99
69) Atrazine	11.069	200	72623	38.039	ng	99
70) Pentachlorophenol	11.198	266	39039	37.863	ng	99
71) Phenanthrene	11.398	178	487215	42.612	ng	100
72) Anthracene	11.451	178	471444	41.869	ng	100
73) Carbazole	11.616	167	437469	39.666	ng	100
74) Di-n-butylphthalate	11.927	149	528355	40.820	ng	99
75) Fluoranthene	12.592	202	514230	39.052	ng	99
77) Benzidine	12.721	184	92303	33.531	ng	99
78) Pyrene	12.827	202	516160	43.035	ng	100
80) Butylbenzylphthalate	13.427	149	175903	39.859	ng	99
81) Benzo(a)anthracene	14.021	228	392086	41.017	ng	99
82) 3,3'-Dichlorobenzidine	13.980	252	125869	43.644	ng	97
83) Chrysene	14.057	228	366916	42.211	ng	99
84) Bis(2-ethylhexyl)phtha...	13.986	149	232385	40.838	ng	98
85) Di-n-octyl phthalate	14.615	149	367492	42.687	ng	98
87) Indeno(1,2,3-cd)pyrene	17.051	276	481277	53.544	ng	99
88) Benzo(b)fluoranthene	15.098	252	364595	36.465	ng	99
89) Benzo(k)fluoranthene	15.127	252	365859m	42.658	ng	
90) Benzo(a)pyrene	15.474	252	337440	43.101	ng	99
91) Dibenzo(a,h)anthracene	17.051	278	402396	54.038	ng	99
92) Benzo(g,h,i)perylene	17.509	276	390342	51.504	ng	99

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

