

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061025\
 Data File : BF142709.D
 Acq On : 10 Jun 2025 13:44
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 10 14:12:31 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 09 18:21:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	71095	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	276013	20.000	ng	0.00
39) Acenaphthene-d10	9.939	164	149668	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	238702	20.000	ng	0.00
76) Chrysene-d12	14.068	240	135085	20.000	ng	0.00
86) Perylene-d12	15.562	264	161985	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	334530	79.895	ng	0.00
7) Phenol-d6	6.522	99	395344	80.514	ng	-0.01
23) Nitrobenzene-d5	7.463	82	399669	79.578	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	125200	74.456	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	869968	79.492	ng	0.00
79) Terphenyl-d14	13.010	244	720168	84.314	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.657	88	67582	42.150	ng	99
3) Pyridine	3.428	79	169546	41.000	ng	99
4) n-Nitrosodimethylamine	3.381	42	80707	40.099	ng	100
6) Aniline	6.557	93	266641	40.447	ng	99
8) 2-Chlorophenol	6.681	128	181660	39.765	ng	99
9) Benzaldehyde	6.445	77	117875	42.583	ng	99
10) Phenol	6.540	94	222155	40.403	ng	99
11) bis(2-Chloroethyl)ether	6.628	93	160191	39.990	ng	99
12) 1,3-Dichlorobenzene	6.834	146	204870	40.044	ng	99
13) 1,4-Dichlorobenzene	6.916	146	210760	40.290	ng	99
14) 1,2-Dichlorobenzene	7.063	146	201071	40.313	ng	99
15) Benzyl Alcohol	7.034	79	148203	41.261	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.169	45	256240	39.283	ng	99
17) 2-Methylphenol	7.145	107	141123	40.287	ng	99
18) Hexachloroethane	7.410	117	73693	41.006	ng	98
19) n-Nitroso-di-n-propyla...	7.304	70	120085	41.720	ng	98
20) 3+4-Methylphenols	7.298	107	178109	41.285	ng	96
22) Acetophenone	7.304	105	238264	39.490	ng	98
24) Nitrobenzene	7.481	77	178378	39.382	ng	99
25) Isophorone	7.716	82	322436	39.571	ng	99
26) 2-Nitrophenol	7.792	139	101800	40.745	ng	96
27) 2,4-Dimethylphenol	7.828	122	169300	39.871	ng	100
28) bis(2-Chloroethoxy)met...	7.928	93	203902	39.897	ng	98
29) 2,4-Dichlorophenol	8.039	162	157567	40.506	ng	99
30) 1,2,4-Trichlorobenzene	8.122	180	171613	40.445	ng	99
31) Naphthalene	8.204	128	534564	39.303	ng	100
32) Benzoic acid	7.939	122	101628	39.468	ng	99
33) 4-Chloroaniline	8.251	127	214232	38.313	ng	99
34) Hexachlorobutadiene	8.316	225	106781	40.861	ng	99
35) Caprolactam	8.616	113	40422	33.832	ng	97
36) 4-Chloro-3-methylphenol	8.728	107	156768	39.070	ng	99
37) 2-Methylnaphthalene	8.892	142	334033	39.031	ng	100
38) 1-Methylnaphthalene	8.992	142	345917	39.032	ng	99
40) 1,2,4,5-Tetrachloroben...	9.057	216	171157	39.571	ng	99
41) Hexachlorocyclopentadiene	9.045	237	108940	42.889	ng	100
43) 2,4,6-Trichlorophenol	9.169	196	114375	40.300	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.210	196	115692	38.165	ng	98
46) 1,1'-Biphenyl	9.357	154	456893	39.258	ng	100
47) 2-Chloronaphthalene	9.380	162	338045	38.766	ng	98
48) 2-Nitroaniline	9.475	65	96497	37.187	ng	94
49) Acenaphthylene	9.798	152	564605	39.067	ng	100
50) Dimethylphthalate	9.657	163	376757	37.415	ng	99
51) 2,6-Dinitrotoluene	9.716	165	82231	37.750	ng	95
52) Acenaphthene	9.969	154	347261	39.278	ng	99
53) 3-Nitroaniline	9.886	138	87885	35.906	ng	98
54) 2,4-Dinitrophenol	9.998	184	46858	38.853	ng	97
55) Dibenzofuran	10.139	168	495828	39.092	ng	99
56) 4-Nitrophenol	10.045	139	60374	35.886	ng	98
57) 2,4-Dinitrotoluene	10.122	165	108516	37.164	ng	100
58) Fluorene	10.486	166	380963	37.579	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.257	232	99666	39.072	ng	99
60) Diethylphthalate	10.351	149	363744	37.208	ng	99
61) 4-Chlorophenyl-phenyle...	10.475	204	185927	38.791	ng	99
62) 4-Nitroaniline	10.498	138	77112	33.211	ng	96
63) Azobenzene	10.633	77	329652	37.644	ng	97
65) 4,6-Dinitro-2-methylph...	10.533	198	59397	39.014	ng	99
66) n-Nitrosodiphenylamine	10.592	169	322521	41.848	ng	100
67) 4-Bromophenyl-phenylether	10.969	248	112692	42.713	ng	98
68) Hexachlorobenzene	11.033	284	123162	40.626	ng	97
69) Atrazine	11.116	200	85223	36.463	ng	98
70) Pentachlorophenol	11.227	266	66476	39.873	ng	99
71) Phenanthrene	11.451	178	501517	39.301	ng	99
72) Anthracene	11.504	178	518482	39.152	ng	100
73) Carbazole	11.657	167	435663	36.534	ng	100
74) Di-n-butylphthalate	11.980	149	472669	37.337	ng	99
75) Fluoranthene	12.639	202	464559	36.334	ng	100
77) Benzidine	12.757	184	205826	44.148	ng	99
78) Pyrene	12.869	202	465611	42.300	ng	99
80) Butylbenzylphthalate	13.480	149	152483	44.983	ng	97
81) Benzo(a)anthracene	14.057	228	341337	38.935	ng	99
82) 3,3'-Dichlorobenzidine	14.021	252	122559	43.734	ng	100
83) Chrysene	14.092	228	331906	39.970	ng	100
84) Bis(2-ethylhexyl)phtha...	14.039	149	239004	61.756	ng	99
85) Di-n-octyl phthalate	14.657	149	461847	61.358	ng	99
87) Indeno(1,2,3-cd)pyrene	17.080	276	517815	45.192	ng	100
88) Benzo(b)fluoranthene	15.121	252	385979	39.883	ng	100
89) Benzo(k)fluoranthene	15.151	252	316516	34.756	ng	100
90) Benzo(a)pyrene	15.498	252	356047	39.261	ng	99
91) Dibenzo(a,h)anthracene	17.098	278	420883	45.630	ng	99
92) Benzo(g,h,i)perylene	17.539	276	421708	44.151	ng	99

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

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