

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070225\
 Data File : BF142971.D
 Acq On : 02 Jul 2025 15:17
 Operator : RC/JU
 Sample : Q2458-04MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-67MSD

Quant Time: Jul 02 15:48:04 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 24 05:28:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	56149	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	216831	20.000	ng	-0.01
39) Acenaphthene-d10	9.916	164	113581	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	197962	20.000	ng	0.00
76) Chrysene-d12	14.056	240	113049	20.000	ng	0.00
86) Perylene-d12	15.545	264	120474	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	252058	74.740	ng	0.00
7) Phenol-d6	6.504	99	304097	76.428	ng	0.00
23) Nitrobenzene-d5	7.433	82	184481	46.667	ng	-0.01
42) 2,4,6-Tribromophenol	10.710	330	95618	81.807	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	416800	48.207	ng	-0.01
79) Terphenyl-d14	12.992	244	381225	46.594	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.645	88	48397	35.879	ng	98
3) Pyridine	3.416	79	130811	38.682	ng	99
4) n-Nitrosodimethylamine	3.363	42	72394	41.399	ng	98
6) Aniline	6.533	93	156171	29.498	ng	# 87
8) 2-Chlorophenol	6.657	128	161846	44.488	ng	98
9) Benzaldehyde	6.422	77	65498	27.747	ng	98
10) Phenol	6.522	94	191654	43.333	ng	93
11) bis(2-Chloroethyl)ether	6.604	93	138878	43.934	ng	98
12) 1,3-Dichlorobenzene	6.816	146	176063	43.274	ng	99
13) 1,4-Dichlorobenzene	6.892	146	177572	43.259	ng	99
14) 1,2-Dichlorobenzene	7.045	146	170609	43.819	ng	100
15) Benzyl Alcohol	7.016	79	129241	44.931	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.145	45	208093	40.974	ng	90
17) 2-Methylphenol	7.128	107	122893	44.577	ng	98
18) Hexachloroethane	7.386	117	60864	42.391	ng	94
19) n-Nitroso-di-n-propyla...	7.286	70	102676	44.369	ng	97
20) 3+4-Methylphenols	7.280	107	155844	45.339	ng	93
22) Acetophenone	7.280	105	213231	44.595	ng	99
24) Nitrobenzene	7.457	77	154127	43.077	ng	98
25) Isophorone	7.698	82	278962	43.991	ng	99
26) 2-Nitrophenol	7.775	139	89684	46.014	ng	98
27) 2,4-Dimethylphenol	7.810	122	148069	43.858	ng	100
28) bis(2-Chloroethoxy)met...	7.904	93	174411	43.261	ng	100
29) 2,4-Dichlorophenol	8.016	162	136833	44.700	ng	99
30) 1,2,4-Trichlorobenzene	8.098	180	146140	43.415	ng	99
31) Naphthalene	8.180	128	460550	43.393	ng	100
32) Benzoic acid	7.927	122	73378	42.643	ng	97
33) 4-Chloroaniline	8.227	127	78716	18.240	ng	98
34) Hexachlorobutadiene	8.292	225	89607	43.520	ng	99
35) Caprolactam	8.604	113	41835m	51.790	ng	
36) 4-Chloro-3-methylphenol	8.716	107	138688	45.586	ng	99
37) 2-Methylnaphthalene	8.874	142	289318	43.969	ng	100
38) 1-Methylnaphthalene	8.974	142	301963	44.331	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	148386	44.276	ng	99
41) Hexachlorocyclopentadiene	9.022	237	155705	81.870	ng	99
43) 2,4,6-Trichlorophenol	9.151	196	96995	44.414	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070225\
 Data File : BF142971.D
 Acq On : 02 Jul 2025 15:17
 Operator : RC/JU
 Sample : Q2458-04MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-67MSD

Quant Time: Jul 02 15:48:04 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 24 05:28:21 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	102673	44.662	ng	99
46) 1,1'-Biphenyl	9.333	154	402287	44.837	ng	100
47) 2-Chloronaphthalene	9.363	162	292943	43.508	ng	99
48) 2-Nitroaniline	9.463	65	87106	46.145	ng	96
49) Acenaphthylene	9.780	152	491743	44.365	ng	99
50) Dimethylphthalate	9.639	163	348094	47.346	ng	100
51) 2,6-Dinitrotoluene	9.704	165	76426	47.334	ng	96
52) Acenaphthene	9.951	154	325254m	48.096	ng	
53) 3-Nitroaniline	9.869	138	48726	27.410	ng	99
54) 2,4-Dinitrophenol	9.986	184	74821	92.547	ng	95
55) Dibenzofuran	10.121	168	439268	45.297	ng	100
56) 4-Nitrophenol	10.039	139	117590	96.622	ng	98
57) 2,4-Dinitrotoluene	10.110	165	105877	49.367	ng	98
58) Fluorene	10.468	166	349636	46.165	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.245	232	85167	45.961	ng	97
60) Diethylphthalate	10.339	149	351584	48.772	ng	99
61) 4-Chlorophenyl-phenyle...	10.457	204	166362	46.033	ng	99
62) 4-Nitroaniline	10.486	138	79394	48.833	ng	99
63) Azobenzene	10.616	77	317643	49.302	ng	98
65) 4,6-Dinitro-2-methylph...	10.521	198	53270	44.488	ng	99
66) n-Nitrosodiphenylamine	10.574	169	300816	43.845	ng	99
67) 4-Bromophenyl-phenylether	10.945	248	100791	44.736	ng	95
68) Hexachlorobenzene	11.016	284	110539	44.155	ng	98
69) Atrazine	11.104	200	93920	53.011	ng	99
70) Pentachlorophenol	11.216	266	105114	84.262	ng	100
71) Phenanthrene	11.433	178	486722	45.388	ng	99
72) Anthracene	11.486	178	494071	44.924	ng	100
73) Carbazole	11.639	167	434982	45.832	ng	100
74) Di-n-butylphthalate	11.963	149	540733	54.691	ng	100
75) Fluoranthene	12.621	202	458695	45.071	ng	100
77) Benzidine	12.745	184	136349	32.102	ng	99
78) Pyrene	12.851	202	450689	42.532	ng	100
80) Butylbenzylphthalate	13.462	149	165587	53.871	ng	99
81) Benzo(a)anthracene	14.045	228	350275	45.929	ng	100
82) 3,3'-Dichlorobenzidine	14.004	252	43306	17.507	ng	# 98
83) Chrysene	14.080	228	308955	45.400	ng	99
84) Bis(2-ethylhexyl)phtha...	14.021	149	216038	48.850	ng	100
85) Di-n-octyl phthalate	14.639	149	333995	40.051	ng	99
87) Indeno(1,2,3-cd)pyrene	17.062	276	364964	39.772	ng	99
88) Benzo(b)fluoranthene	15.109	252	330057	47.346	ng	99
89) Benzo(k)fluoranthene	15.139	252	312488	47.913	ng	99
90) Benzo(a)pyrene	15.486	252	311232	46.214	ng	99
91) Dibenzo(a,h)anthracene	17.074	278	294876	39.549	ng	99
92) Benzo(g,h,i)perylene	17.521	276	286817	38.577	ng	99

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

