

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF033125\
 Data File : BF142177.D
 Acq On : 31 Mar 2025 13:22
 Operator : RC/JU
 Sample : PB167369BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB167369BS

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 04/01/2025
 Supervised By :Jagrut Upadhyay 04/01/2025

Quant Time: Mar 31 13:49:11 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 10 15:46:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	120659	20.000	ng	-0.01
21) Naphthalene-d8	8.151	136	484067	20.000	ng	-0.01
39) Acenaphthene-d10	9.904	164	285734	20.000	ng	-0.01
64) Phenanthrene-d10	11.392	188	517636	20.000	ng	-0.01
76) Chrysene-d12	14.039	240	319787	20.000	ng	0.00
86) Perylene-d12	15.509	264	347760	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	1036795	143.410	ng	0.00
7) Phenol-d6	6.498	99	1271824	138.169	ng	0.00
23) Nitrobenzene-d5	7.428	82	834922	97.065	ng	-0.02
42) 2,4,6-Tribromophenol	10.698	330	627855	173.206	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1857788	98.880	ng	-0.01
79) Terphenyl-d14	12.980	244	2425298	112.127	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.704	88	126082	41.622	ng	96
3) Pyridine	3.463	79	339519	45.693	ng	96
4) n-Nitrosodimethylamine	3.410	42	159631	45.068	ng	93
6) Aniline	6.528	93	246191	27.112	ng	# 89
8) 2-Chlorophenol	6.651	128	421298	52.543	ng	97
9) Benzaldehyde	6.416	77	253399	49.222	ng	99
10) Phenol	6.510	94	476524	49.208	ng	98
11) bis(2-Chloroethyl)ether	6.604	93	356801	49.069	ng	96
12) 1,3-Dichlorobenzene	6.810	146	446871	51.623	ng	98
13) 1,4-Dichlorobenzene	6.886	146	452389	51.651	ng	99
14) 1,2-Dichlorobenzene	7.033	146	432910	52.476	ng	99
15) Benzyl Alcohol	7.004	79	362663	48.734	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.139	45	380763	43.374	ng	95
17) 2-Methylphenol	7.116	107	326040	51.141	ng	99
18) Hexachloroethane	7.380	117	165135	50.279	ng	93
19) n-Nitroso-di-n-propyla...	7.281	70	272051	45.996	ng	99
20) 3+4-Methylphenols	7.269	107	414507	50.761	ng	# 87
22) Acetophenone	7.275	105	553030	47.707	ng	96
24) Nitrobenzene	7.451	77	414348	48.456	ng	97
25) Isophorone	7.686	82	775507	51.004	ng	100
26) 2-Nitrophenol	7.763	139	227763	54.270	ng	98
27) 2,4-Dimethylphenol	7.798	122	389738	67.441	ng	99
28) bis(2-Chloroethoxy)met...	7.898	93	464694	48.727	ng	99
29) 2,4-Dichlorophenol	8.004	162	369700	51.540	ng	99
30) 1,2,4-Trichlorobenzene	8.092	180	409755	51.836	ng	98
31) Naphthalene	8.169	128	1202192	48.347	ng	100
32) Benzoic acid	7.922	122	268561m	52.868	ng	
33) 4-Chloroaniline	8.222	127	85644	9.757	ng	98
34) Hexachlorobutadiene	8.286	225	272000	52.763	ng	100
35) Caprolactam	8.592	113	113738m	52.009	ng	
36) 4-Chloro-3-methylphenol	8.704	107	409653	51.197	ng	97
37) 2-Methylnaphthalene	8.863	142	754830	45.416	ng	100
38) 1-Methylnaphthalene	8.963	142	790773	49.304	ng	99
40) 1,2,4,5-Tetrachloroben...	9.027	216	444807	51.130	ng	98
41) Hexachlorocyclopentadiene	9.016	237	622980	182.981	ng	99
43) 2,4,6-Trichlorophenol	9.139	196	298031	52.420	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.180	196	307687	53.434	ng	99
46) 1,1'-Biphenyl	9.327	154	1076545	49.586	ng	99
47) 2-Chloronaphthalene	9.351	162	825118	50.990	ng	98
48) 2-Nitroaniline	9.451	65	227855	48.632	ng	95
49) Acenaphthylene	9.769	152	1301383	54.194	ng	100
50) Dimethylphthalate	9.627	163	1024272	51.190	ng	100
51) 2,6-Dinitrotoluene	9.692	165	223399	53.623	ng	92
52) Acenaphthene	9.939	154	829966	49.182	ng	100
53) 3-Nitroaniline	9.857	138	60833	14.395	ng	97
54) 2,4-Dinitrophenol	9.969	184	263709	108.826	ng	# 45
55) Dibenzofuran	10.116	168	1182589	49.044	ng	98
56) 4-Nitrophenol	10.027	139	336591	105.616	ng	95
57) 2,4-Dinitrotoluene	10.098	165	314970	57.884	ng	98
58) Fluorene	10.457	166	960135	51.633	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.233	232	282918	53.994	ng	96
60) Diethylphthalate	10.327	149	1023786	51.313	ng	100
61) 4-Chlorophenyl-phenyle...	10.445	204	511854	52.797	ng	98
62) 4-Nitroaniline	10.474	138	206720	50.245	ng	98
63) Azobenzene	10.610	77	878646	47.368	ng	96
65) 4,6-Dinitro-2-methylph...	10.504	198	183397	59.432	ng	97
66) n-Nitrosodiphenylamine	10.569	169	841785	50.253	ng	100
67) 4-Bromophenyl-phenylether	10.939	248	340379	52.359	ng	96
68) Hexachlorobenzene	11.004	284	390998	54.838	ng	95
69) Atrazine	11.092	200	323666	63.029	ng	99
70) Pentachlorophenol	11.198	266	455852	103.768	ng	100
71) Phenanthrene	11.421	178	1445434	51.698	ng	99
72) Anthracene	11.474	178	1476065	52.622	ng	100
73) Carbazole	11.627	167	1224318	50.610	ng	100
74) Di-n-butylphthalate	11.951	149	1611821	50.215	ng	100
75) Fluoranthene	12.610	202	1509227	50.887	ng	99
77) Benzidine	12.727	184	383514	85.958	ng	100
78) Pyrene	12.839	202	1509100	54.555	ng	100
80) Butylbenzylphthalate	13.451	149	547547	50.573	ng	99
81) Benzo(a)anthracene	14.027	228	1118309	53.292	ng	100
82) 3,3'-Dichlorobenzidine	13.986	252	58588	9.661	ng	99
83) Chrysene	14.062	228	1007817	52.982	ng	100
84) Bis(2-ethylhexyl)phtha...	14.009	149	720391	48.258	ng	98
85) Di-n-octyl phthalate	14.627	149	1085911	52.280	ng	97
87) Indeno(1,2,3-cd)pyrene	17.009	276	1167687	51.906	ng	97
88) Benzo(b)fluoranthene	15.080	252	1139652	47.816	ng	99
89) Benzo(k)fluoranthene	15.109	252	895900	43.924	ng	100
90) Benzo(a)pyrene	15.451	252	1058724	56.770	ng	100
91) Dibenzo(a,h)anthracene	17.027	278	952603	51.323	ng	99
92) Benzo(g,h,i)perylene	17.462	276	918852	50.095	ng	98

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

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