

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112625\
 Data File : BF144378.D
 Acq On : 26 Nov 2025 23:44
 Operator : RC/JU
 Sample : Q3721-01MS 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-2-112525MS

Quant Time: Nov 27 00:55:21 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 05 18:37:33 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	62340	20.000	ng	-0.01
21) Naphthalene-d8	8.145	136	204401	20.000	ng	-0.01
39) Acenaphthene-d10	9.904	164	101875	20.000	ng	-0.01
64) Phenanthrene-d10	11.392	188	181460	20.000	ng	-0.01
76) Chrysene-d12	14.045	240	168757	20.000	ng	0.00
86) Perylene-d12	15.527	264	133172	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	71054	17.836	ng	0.00
7) Phenol-d6	6.493	99	76483	16.944	ng	0.00
23) Nitrobenzene-d5	7.422	82	43555	12.307	ng	-0.01
42) 2,4,6-Tribromophenol	10.692	330	20084	15.710	ng	-0.01
45) 2-Fluorobiphenyl	9.216	172	85255	12.360	ng	-0.01
79) Terphenyl-d14	12.980	244	109336	10.291	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.593	88	12601	7.651	ng	# 95
3) Pyridine	3.387	79	32040	6.973	ng	95
4) n-Nitrosodimethylamine	3.304	42	17488	7.286	ng	93
6) Aniline	6.522	93	36984	5.751	ng	# 82
8) 2-Chlorophenol	6.651	128	32682	8.365	ng	96
9) Benzaldehyde	6.416	77	16822	6.600	ng	# 82
10) Phenol	6.510	94	46193	8.376	ng	98
11) bis(2-Chloroethyl)ether	6.593	93	33901	8.436	ng	98
12) 1,3-Dichlorobenzene	6.804	146	41027	8.510	ng	94
13) 1,4-Dichlorobenzene	6.881	146	39899	8.261	ng	96
14) 1,2-Dichlorobenzene	7.034	146	38372	8.289	ng	97
15) Benzyl Alcohol	7.004	79	26866	7.517	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.134	45	57715	7.799	ng	81
17) 2-Methylphenol	7.116	107	25640	8.250	ng	99
18) Hexachloroethane	7.375	117	16971	10.577	ng	93
19) n-Nitroso-di-n-propyla...	7.263	70	27465	9.243	ng	97
20) 3+4-Methylphenols	7.275	107	31457	8.586	ng	97
22) Acetophenone	7.269	105	59044	13.465	ng	94
24) Nitrobenzene	7.439	77	32995	8.587	ng	97
25) Isophorone	7.681	82	60960	9.106	ng	# 93
26) 2-Nitrophenol	7.763	139	16377	8.610	ng	98
27) 2,4-Dimethylphenol	7.798	122	28810	10.105	ng	99
28) bis(2-Chloroethoxy)met...	7.887	93	37895	9.065	ng	# 90
29) 2,4-Dichlorophenol	8.010	162	26740	8.138	ng	97
30) 1,2,4-Trichlorobenzene	8.087	180	28998	8.594	ng	95
31) Naphthalene	8.163	128	90701	9.328	ng	98
32) Benzoic acid	7.892	122	11140	5.331	ng	# 59
33) 4-Chloroaniline	8.222	127	21592	6.089	ng	97
34) Hexachlorobutadiene	8.281	225	19916	7.832	ng	98
35) Caprolactam	8.557	113	11431	12.697	ng	# 28
36) 4-Chloro-3-methylphenol	8.704	107	22877	7.768	ng	98
37) 2-Methylnaphthalene	8.857	142	57344	8.821	ng	98
38) 1-Methylnaphthalene	8.957	142	55980	8.925	ng	99
40) 1,2,4,5-Tetrachloroben...	9.022	216	27554	8.253	ng	98
41) Hexachlorocyclopentadiene	9.010	237	489	4.761	ng	94
43) 2,4,6-Trichlorophenol	9.139	196	18190	8.005	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	19333	7.894	ng	# 83
46) 1,1'-Biphenyl	9.322	154	70738	9.947	ng	98
47) 2-Chloronaphthalene	9.345	162	51087	8.422	ng	99
48) 2-Nitroaniline	9.445	65	14654	8.373	ng	95
49) Acenaphthylene	9.763	152	81412	9.849	ng	99
50) Dimethylphthalate	9.616	163	59631	8.836	ng	# 97
51) 2,6-Dinitrotoluene	9.686	165	11997	8.781	ng	89
52) Acenaphthene	9.933	154	42726	7.729	ng	99
53) 3-Nitroaniline	9.857	138	12107	8.107	ng	# 94
55) Dibenzofuran	10.110	168	66970	8.650	ng	97
56) 4-Nitrophenol	10.039	139	13801	12.775	ng	92
57) 2,4-Dinitrotoluene	10.092	165	14402	7.874	ng	# 98
58) Fluorene	10.451	166	55760	9.473	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.233	232	13490	7.785	ng	# 79
60) Diethylphthalate	10.316	149	54583	8.773	ng	100
61) 4-Chlorophenyl-phenyle...	10.439	204	28179	8.404	ng	94
62) 4-Nitroaniline	10.469	138	12335	8.673	ng	92
63) Azobenzene	10.604	77	48839	8.435	ng	88
66) n-Nitrosodiphenylamine	10.557	169	53940	9.242	ng	96
67) 4-Bromophenyl-phenylether	10.933	248	17913	7.734	ng	96
68) Hexachlorobenzene	11.004	284	20566	7.538	ng	97
69) Atrazine	11.086	200	17128	9.228	ng	96
70) Pentachlorophenol	11.204	266	20060	14.766	ng	90
71) Phenanthrene	11.416	178	93078	10.302	ng	99
72) Anthracene	11.469	178	87986	9.578	ng	100
73) Carbazole	11.627	167	78560	10.057	ng	98
74) Di-n-butylphthalate	11.951	149	91726	9.713	ng	99
75) Fluoranthene	12.610	202	107938	11.565	ng	98
77) Benzidine	12.733	184	41547	8.323	ng	97
78) Pyrene	12.839	202	122440	8.999	ng	99
80) Butylbenzylphthalate	13.457	149	44501	9.436	ng	95
81) Benzo(a)anthracene	14.033	228	105655	9.033	ng	99
82) 3,3'-Dichlorobenzidine	13.998	252	24595	6.135	ng	# 99
83) Chrysene	14.068	228	98150	9.090	ng	96
84) Bis(2-ethylhexyl)phtha...	14.016	149	65427	10.135	ng	99
85) Di-n-octyl phthalate	14.627	149	101594	9.121	ng	# 88
87) Indeno(1,2,3-cd)pyrene	17.027	276	67213	7.031	ng	# 95
88) Benzo(b)fluoranthene	15.092	252	76308	10.079	ng	# 98
89) Benzo(k)fluoranthene	15.121	252	69911	9.865	ng	# 95
90) Benzo(a)pyrene	15.463	252	67374	9.646	ng	# 96
91) Dibenzo(a,h)anthracene	17.039	278	54023	7.037	ng	# 95
92) Benzo(g,h,i)perylene	17.480	276	53702	7.083	ng	# 94

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

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