

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111324\
 Data File : BF140345.D
 Acq On : 13 Nov 2024 16:40
 Operator : RC/JU
 Sample : P3845-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-BN-WP

Quant Time: Nov 13 17:19:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 14:40:06 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	135927	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	521483	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	291001	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	534019	20.000	ng	0.00
76) Chrysene-d12	14.080	240	320922	20.000	ng	0.03
86) Perylene-d12	15.621	264	276076	20.000	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	1097292	137.831	ng	0.00
7) Phenol-d6	6.510	99	1533295	142.156	ng	0.00
23) Nitrobenzene-d5	7.439	82	995732	99.486	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	411701	142.271	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	1728492	95.485	ng	0.00
79) Terphenyl-d14	12.992	244	2021563	109.342	ng	0.00
Target Compounds						
						Qvalue
4) n-Nitrosodimethylamine	3.428	42	518523	117.613	ng	98
11) bis(2-Chloroethyl)ether	6.616	93	945580	109.713	ng	98
12) 1,3-Dichlorobenzene	6.822	146	1155625	122.476	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.151	45	837740	70.371	ng	99
18) Hexachloroethane	7.392	117	398527	112.675	ng	100
19) n-Nitroso-di-n-propyla...	7.287	70	753459	115.662	ng	99
24) Nitrobenzene	7.457	77	225776	21.666	ng	99
31) Naphthalene	8.181	128	3185954	120.434	ng	98
34) Hexachlorobutadiene	8.292	225	353570	68.123	ng	99
37) 2-Methylnaphthalene	8.869	142	1370865	80.817	ng	100
38) 1-Methylnaphthalene	8.975	142	2170313	130.658	ng	100
41) Hexachlorocyclopentadiene	9.022	237	723777	350.306	ng	99
49) Acenaphthylene	9.780	152	3065708	125.646	ng	98
51) 2,6-Dinitrotoluene	9.698	165	367918	85.084	ng	96
52) Acenaphthene	9.945	154	328135	19.911	ng	100
53) 3-Nitroaniline	9.863	138	27688	6.097	ng	# 97
55) Dibenzofuran	10.122	168	1283563	55.019	ng	100
57) 2,4-Dinitrotoluene	10.104	165	367514	64.578	ng	99
58) Fluorene	10.475	166	2576051	139.776	ng	99
60) Diethylphthalate	10.339	149	1561064	80.486	ng	100
61) 4-Chlorophenyl-phenyle...	10.457	204	324466	35.661	ng	99
66) n-Nitrosodiphenylamine	10.575	169	1261321	81.747	ng	99
68) Hexachlorobenzene	11.016	284	771805	128.503	ng	98
72) Anthracene	11.480	178	2289055	93.105	ng	99
78) Pyrene	12.845	202	930946	33.913	ng	100
83) Chrysene	14.110	228	2126934	110.522	ng	99
84) Bis(2-ethylhexyl)phtha...	14.057	149	1277183	90.738	ng	98
85) Di-n-octyl phthalate	14.721	149	2318732	116.967	ng	100
87) Indeno(1,2,3-cd)pyrene	17.151	276	1861642	109.162	ng	100
88) Benzo(b)fluoranthene	15.180	252	771191	42.703	ng	99
89) Benzo(k)fluoranthene	15.215	252	1149520	77.916	ng	99
90) Benzo(a)pyrene	15.557	252	918715	65.508	ng	99
91) Dibenzo(a,h)anthracene	17.168	278	1351813	95.896	ng	99
92) Benzo(g,h,i)perylene	17.603	276	1175181	81.545	ng	99

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

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