

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100422\
 Data File : BN022008.D
 Acq On : 04 Oct 2022 15:37
 Operator : CG/JU
 Sample : N4971-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWOW-BR-02-231-251-100322-FD

Quant Time: Oct 04 23:04:59 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 04 01:32:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.985	152	4458	0.400	ng	0.00
7) Naphthalene-d8	10.816	136	14108	0.400	ng	0.00
13) Acenaphthene-d10	14.646	164	7768	0.400	ng	0.00
19) Phenanthrene-d10	17.403	188	15740	0.400	ng	# 0.00
29) Chrysene-d12	21.608	240	13700	0.400	ng	0.00
35) Perylene-d12	24.084	264	11463	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.515	112	1460	0.141	ng	0.00
5) Phenol-d6	7.140	99	1105	0.082	ng	0.00
8) Nitrobenzene-d5	9.161	82	3890	0.343	ng	0.00
11) 2-Methylnaphthalene-d10	12.406	152	7009	0.270	ng	0.00
14) 2,4,6-Tribromophenol	16.150	330	1267	0.428	ng	0.00
15) 2-Fluorobiphenyl	13.278	172	11227	0.331	ng	0.00
27) Fluoranthene-d10	19.441	212	17382	0.415	ng	0.00
31) Terphenyl-d14	20.032	244	11324	0.411	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.859	128	5799	0.136	ng	97
16) Acenaphthylene	14.368	152	1316	0.036	ng	97
18) Fluorene	15.694	166	1159	0.037	ng	# 99
24) Pentachlorophenol	17.056	266	82	0.022	ng	86
25) Phenanthrene	17.441	178	5189	0.096	ng	98
26) Anthracene	17.540	178	1661	0.038	ng	# 90
28) Fluoranthene	19.470	202	3076	0.053	ng	# 86
30) Pyrene	19.833	202	2636	0.044	ng	# 93
32) Benzo(a)anthracene	21.591	228	2044	0.040	ng	# 83
33) Chrysene	21.644	228	1757	0.031	ng	# 83
34) Bis(2-ethylhexyl)phtha...	21.501	149	33190	1.343	ng	98
36) Indeno(1,2,3-cd)pyrene	26.686	276	1165	0.020	ng	# 88
37) Benzo(b)fluoranthene	23.324	252	1858	0.034	ng	# 1
38) Benzo(k)fluoranthene	23.367	252	1226	0.023	ng	# 1
39) Benzo(a)pyrene	23.970	252	1382	0.033	ng	# 1
41) Benzo(g,h,i)perylene	27.496	276	1103	0.022	ng	# 60

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

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