

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100422\
 Data File : BN022013.D
 Acq On : 04 Oct 2022 19:40
 Operator : CG/JU
 Sample : SP6026
 Misc : 8270SIM SURROGATES
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP6026

Quant Time: Oct 04 23:06:42 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	5868	0.400	ng	0.00
7) Naphthalene-d8	10.816	136	17404	0.400	ng	0.00
13) Acenaphthene-d10	14.653	164	8847	0.400	ng	0.00
19) Phenanthrene-d10	17.412	188	18078	0.400	ng	0.00
29) Chrysene-d12	21.609	240	12062	0.400	ng	0.00
35) Perylene-d12	24.087	264	9722	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.515	112	6221	0.458	ng	0.00
5) Phenol-d6	7.148	99	7513	0.426	ng	0.00
8) Nitrobenzene-d5	9.161	82	4703	0.336	ng	0.00
11) 2-Methylnaphthalene-d10	12.410	152	10461	0.327	ng	0.00
14) 2,4,6-Tribromophenol	16.158	330	1152	0.341	ng	0.01
15) 2-Fluorobiphenyl	13.285	172	13543	0.350	ng	0.00
27) Fluoranthene-d10	19.445	212	15294	0.318	ng	0.00
31) Terphenyl-d14	20.033	244	10189	0.420	ng	0.00

Target Compounds Qvalue

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

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