

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110222\
 Data File : BN022459.D
 Acq On : 02 Nov 2022 10:16
 Operator : CG/JU
 Sample : PB148513BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB148513BL

Quant Time: Nov 02 11:32:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 02 05:31:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	3586	0.400	ng	0.00
7) Naphthalene-d8	10.752	136	11599	0.400	ng	0.00
13) Acenaphthene-d10	14.589	164	5079	0.400	ng	0.00
19) Phenanthrene-d10	17.350	188	9385	0.400	ng	0.00
29) Chrysene-d12	21.555	240	8840	0.400	ng	0.00
35) Perylene-d12	23.999	264	9073	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.458	112	4048	0.424	ng	0.00
5) Phenol-d6	7.090	99	4637	0.387	ng	0.00
8) Nitrobenzene-d5	9.108	82	3428	0.342	ng	0.00
11) 2-Methylnaphthalene-d10	12.349	152	5190	0.285	ng	0.00
14) 2,4,6-Tribromophenol	16.109	330	538	0.246	ng	0.01
15) 2-Fluorobiphenyl	13.221	172	7343	0.338	ng	0.00
27) Fluoranthene-d10	19.390	212	8732	0.337	ng	0.00
31) Terphenyl-d14	19.985	244	8369	0.363	ng	0.00

Target Compounds Qvalue

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

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