

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012925\
 Data File : BN036113.D
 Acq On : 29 Jan 2025 18:42
 Operator : RC/JU
 Sample : PB166297BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB166297BL

Quant Time: Jan 30 00:35:33 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN012225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 23 00:34:56 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.803	152	2136	0.400	ng	-0.01
7) Naphthalene-d8	10.600	136	4915	0.400	ng	-0.01
13) Acenaphthene-d10	14.442	164	2707	0.400	ng	0.00
19) Phenanthrene-d10	17.186	188	5606	0.400	ng	0.00
29) Chrysene-d12	21.376	240	4293	0.400	ng	0.00
35) Perylene-d12	23.678	264	4681	0.400	ng	# 0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.419	112	2354	0.424	ng	0.03
5) Phenol-d6	7.023	99	2714	0.416	ng	0.05
8) Nitrobenzene-d5	8.956	82	1959	0.422	ng	0.00
11) 2-Methylnaphthalene-d10	12.192	152	2885	0.432	ng	0.00
14) 2,4,6-Tribromophenol	15.945	330	538	0.310	ng	0.02
15) 2-Fluorobiphenyl	13.073	172	4491	0.372	ng	0.00
27) Fluoranthene-d10	19.220	212	6282	0.433	ng	0.00
31) Terphenyl-d14	19.820	244	5035	0.565	ng	0.00

Target Compounds Qvalue

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

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