

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060425\
 Data File : BN037167.D
 Acq On : 04 Jun 2025 11:03
 Operator : RC/JU
 Sample : PB168238BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB168238BL

Quant Time: Jun 04 12:25:23 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN060325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 04 01:52:03 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.589	152	2941	0.400	ng	0.00
7) Naphthalene-d8	10.372	136	7107	0.400	ng	0.00
13) Acenaphthene-d10	14.245	164	3690	0.400	ng	0.01
19) Phenanthrene-d10	16.984	188	6730	0.400	ng	0.00
29) Chrysene-d12	21.189	240	4435	0.400	ng	0.00
35) Perylene-d12	23.374	264	4132	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.192	112	2715	0.373	ng	0.00
5) Phenol-d6	6.773	99	3089	0.350	ng	0.00
8) Nitrobenzene-d5	8.739	82	2547	0.340	ng	0.00
11) 2-Methylnaphthalene-d10	11.971	152	3335	0.337	ng	0.00
14) 2,4,6-Tribromophenol	15.743	330	378	0.254	ng	0.00
15) 2-Fluorobiphenyl	12.863	172	5709	0.363	ng	0.00
27) Fluoranthene-d10	19.026	212	5729	0.335	ng	0.00
31) Terphenyl-d14	19.630	244	3861	0.370	ng	0.00

Target Compounds	Qvalue

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

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