

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030625\
 Data File : BN036535.D
 Acq On : 06 Mar 2025 14:48
 Operator : RC/JU
 Sample : PB167006BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB167006BL

Quant Time: Mar 06 15:34:58 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN021025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 05 16:08:57 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.732	152	1970	0.400	ng	0.00
7) Naphthalene-d8	10.530	136	4027	0.400	ng	0.02
13) Acenaphthene-d10	14.377	164	2414	0.400	ng	0.01
19) Phenanthrene-d10	17.124	188	4679	0.400	ng	# 0.01
29) Chrysene-d12	21.304	240	4156	0.400	ng	# 0.00
35) Perylene-d12	23.569	264	3735	0.400	ng	# 0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.319	112	1966	0.422	ng	0.00
5) Phenol-d6	6.908	99	1962	0.359	ng	0.00
8) Nitrobenzene-d5	8.886	82	1628	0.410	ng	0.01
11) 2-Methylnaphthalene-d10	12.131	152	2309	0.373	ng	0.03
14) 2,4,6-Tribromophenol	15.882	330	337	0.282	ng	0.02
15) 2-Fluorobiphenyl	13.008	172	3433	0.378	ng	0.02
27) Fluoranthene-d10	19.150	212	5795	0.445	ng	0.00
31) Terphenyl-d14	19.754	244	3979	0.449	ng	0.00

Target Compounds Qvalue

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

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