

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021225\
 Data File : BN036438.D
 Acq On : 12 Feb 2025 13:19
 Operator : RC/JU
 Sample : PB166297BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB166297BL

Quant Time: Feb 12 14:00:19 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN021025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 11 01:17:14 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.753	152	2043	0.400	ng	0.00
7) Naphthalene-d8	10.562	136	4466	0.400	ng	# 0.02
13) Acenaphthene-d10	14.398	164	2771	0.400	ng	0.01
19) Phenanthrene-d10	17.148	188	6288	0.400	ng	0.01
29) Chrysene-d12	21.322	240	5467	0.400	ng	0.00
35) Perylene-d12	23.592	264	5274	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.348	112	1897	0.393	ng	0.00
5) Phenol-d6	6.937	99	1876	0.331	ng	0.00
8) Nitrobenzene-d5	8.929	82	1415	0.321	ng	0.02
11) 2-Methylnaphthalene-d10	12.162	152	2276	0.332	ng	0.02
14) 2,4,6-Tribromophenol	15.907	330	291	0.212	ng	0.02
15) 2-Fluorobiphenyl	13.040	172	3273	0.314	ng	0.02
27) Fluoranthene-d10	19.169	212	6544	0.374	ng	0.00
31) Terphenyl-d14	19.768	244	4993	0.428	ng	0.00

Target Compounds Qvalue

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

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