

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110122\
 Data File : BN022445.D
 Acq On : 01 Nov 2022 22:20
 Operator : CG/JU
 Sample : PB148513BL
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB148513BL

Quant Time: Nov 02 05:38:21 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	3692	0.400	ng	0.00
7) Naphthalene-d8	10.752	136	10949	0.400	ng	# 0.00
13) Acenaphthene-d10	14.600	164	5602	0.400	ng	0.01
19) Phenanthrene-d10	17.362	188	10181	0.400	ng	# 0.01
29) Chrysene-d12	21.555	240	5932	0.400	ng	0.00
35) Perylene-d12	24.002	264	4944	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.465	112	3461	0.352	ng	0.00
5) Phenol-d6	7.097	99	3948	0.320	ng	0.00
8) Nitrobenzene-d5	9.118	82	2832	0.299	ng	0.01
11) 2-Methylnaphthalene-d10	12.357	152	5186	0.302	ng	0.00
14) 2,4,6-Tribromophenol	16.121	330	407	0.169	ng	0.02
15) 2-Fluorobiphenyl	13.231	172	7967	0.333	ng	0.01
27) Fluoranthene-d10	19.394	212	7527	0.268	ng	0.00
31) Terphenyl-d14	19.989	244	6910	0.447	ng	0.00

Target Compounds Qvalue

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

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