

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090524\
 Data File : BN033761.D
 Acq On : 05 Sep 2024 10:32
 Operator : MA/JU
 Sample : PB163111BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB163111BL

Quant Time: Sep 05 11:28:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN090424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 18:36:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.516	152	5687	0.400	ng	0.00
7) Naphthalene-d8	10.271	136	14388	0.400	ng	0.00
13) Acenaphthene-d10	14.156	164	6570	0.400	ng	0.00
19) Phenanthrene-d10	16.904	188	12633	0.400	ng	0.00
29) Chrysene-d12	21.112	240	9448	0.400	ng	0.00
35) Perylene-d12	23.268	264	9984	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.161	112	5779	0.332	ng	0.00
5) Phenol-d6	6.707	99	6524	0.317	ng	0.00
8) Nitrobenzene-d5	8.649	82	4809	0.393	ng	0.00
11) 2-Methylnaphthalene-d10	11.873	152	7879	0.387	ng	0.00
14) 2,4,6-Tribromophenol	15.651	330	738	0.222	ng	0.00
15) 2-Fluorobiphenyl	12.767	172	11000	0.402	ng	0.00
27) Fluoranthene-d10	18.947	212	12845	0.412	ng	0.00
31) Terphenyl-d14	19.560	244	8982	0.446	ng	0.00

Target Compounds Qvalue

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

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