

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
 Data File : BN034487.D
 Acq On : 09 Oct 2024 03:08
 Operator : RC/JU
 Sample : P4226-12
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT119S1-20240925

Quant Time: Oct 09 04:07:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 26 14:44:02 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.380	152	5548	0.400	ng	-0.01
7) Naphthalene-d8	10.137	136	15746	0.400	ng	#-0.01
13) Acenaphthene-d10	14.030	164	8556	0.400	ng	-0.02
19) Phenanthrene-d10	16.796	188	13947	0.400	ng	0.00
29) Chrysene-d12	21.008	240	8936	0.400	ng	0.00
35) Perylene-d12	23.119	264	12144	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.026	112	2152	0.137	ng	0.00
5) Phenol-d6	6.593	99	861	0.047	ng	0.00
8) Nitrobenzene-d5	8.536	82	3049	0.253	ng	0.00
11) 2-Methylnaphthalene-d10	11.742	152	6813	0.293	ng	-0.01
14) 2,4,6-Tribromophenol	15.542	330	1165	0.301	ng	0.00
15) 2-Fluorobiphenyl	12.651	172	9375	0.269	ng	0.00
27) Fluoranthene-d10	18.829	212	15459	0.439	ng	-0.02
31) Terphenyl-d14	19.452	244	10607	0.594	ng	-0.02

Target Compounds Qvalue

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

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