

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
 Data File : BN034492.D
 Acq On : 09 Oct 2024 06:09
 Operator : RC/JU
 Sample : P4226-17
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW9-MW-01D2-20240925

Quant Time: Oct 09 06:57:44 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	7.373	152	4487	0.400	ng	-0.02
7) Naphthalene-d8	10.137	136	11504	0.400	ng	-0.01
13) Acenaphthene-d10	14.026	164	6323	0.400	ng	-0.02
19) Phenanthrene-d10	16.791	188	10885	0.400	ng	#-0.01
29) Chrysene-d12	21.015	240	9274	0.400	ng	0.00
35) Perylene-d12	23.119	264	11019	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.026	112	1890	0.148	ng	0.00
5) Phenol-d6	6.586	99	798	0.054	ng	-0.01
8) Nitrobenzene-d5	8.536	82	2336	0.265	ng	0.00
11) 2-Methylnaphthalene-d10	11.742	152	5195	0.306	ng	-0.01
14) 2,4,6-Tribromophenol	15.549	330	876	0.306	ng	0.00
15) 2-Fluorobiphenyl	12.647	172	7475	0.291	ng	-0.01
27) Fluoranthene-d10	18.832	212	13083	0.476	ng	-0.01
31) Terphenyl-d14	19.450	244	9173	0.495	ng	-0.02

Target Compounds Qvalue

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

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