

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101124\
 Data File : BN034554.D
 Acq On : 11 Oct 2024 14:35
 Operator : RC/JU
 Sample : P4267-01RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BPOW6-11-HYD-20241001

Quant Time: Oct 11 15:42:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 00:21:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.366	152	4037	0.400	ng	0.00
7) Naphthalene-d8	10.127	136	10825	0.400	ng	0.00
13) Acenaphthene-d10	14.019	164	6085	0.400	ng	0.00
19) Phenanthrene-d10	16.784	188	10936	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	4789	0.400	ng	# 0.00
35) Perylene-d12	23.107	264	8624	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.011	112	1342	0.117	ng	0.00
5) Phenol-d6	6.586	99	10	0.001	ng	0.01
8) Nitrobenzene-d5	8.525	82	2183	0.264	ng	0.01
11) 2-Methylnaphthalene-d10	11.742	152	1027	0.064	ng	0.02
14) 2,4,6-Tribromophenol	15.542	330	778	0.282	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	8114	0.328	ng	0.00
27) Fluoranthene-d10	18.825	212	10890	0.395	ng	0.00
31) Terphenyl-d14	19.443	244	9393	0.982	ng	0.00
Target Compounds						
6) bis(2-Chloroethyl)ether	6.781	93	24775	2.101	ng	# 38

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

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