

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101424\
 Data File : BN034585.D
 Acq On : 14 Oct 2024 23:31
 Operator : RC/JU
 Sample : P4267-01RE
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

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

Quant Time: Oct 15 00:07:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN101424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 14 17:52:02 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.366	152	4163	0.400	ng	0.00
7) Naphthalene-d8	10.127	136	11428	0.400	ng	0.01
13) Acenaphthene-d10	14.015	164	6745	0.400	ng	0.00
19) Phenanthrene-d10	16.778	188	11809	0.400	ng	# 0.00
29) Chrysene-d12	21.006	240	4859	0.400	ng	0.00
35) Perylene-d12	23.105	264	7116	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.011	112	1345	0.119	ng	0.00
5) Phenol-d6	6.557	99	3	0.000	ng	0.00
8) Nitrobenzene-d5	8.525	82	2524	0.339	ng	0.01
11) 2-Methylnaphthalene-d10	11.742	152	1206	0.070	ng	0.02
14) 2,4,6-Tribromophenol	15.537	330	743	0.249	ng	0.00
15) 2-Fluorobiphenyl	12.636	172	8477	0.364	ng	0.00
27) Fluoranthene-d10	18.822	212	11564	0.403	ng	0.00
31) Terphenyl-d14	19.445	244	9811	0.473	ng	0.00

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

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