

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034526.D
 Acq On : 10 Oct 2024 18:07
 Operator : RC/JU
 Sample : P4263-08
 Misc :
 ALS Vial : 102 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP11-20240930

Quant Time: Oct 10 18:38:24 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	4598	0.400	ng	-0.02
7) Naphthalene-d8	10.137	136	12965	0.400	ng	-0.01
13) Acenaphthene-d10	14.019	164	7649	0.400	ng	-0.03
19) Phenanthrene-d10	16.784	188	14591	0.400	ng	#-0.02
29) Chrysene-d12	21.008	240	8330	0.400	ng	0.00
35) Perylene-d12	23.116	264	10743	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	2211	0.169	ng	-0.01
5) Phenol-d6	6.593	99	950	0.063	ng	0.00
8) Nitrobenzene-d5	8.536	82	2135	0.215	ng	0.00
11) 2-Methylnaphthalene-d10	11.735	152	5763	0.301	ng	-0.02
14) 2,4,6-Tribromophenol	15.542	330	973	0.281	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	8194	0.263	ng	-0.02
27) Fluoranthene-d10	18.829	212	14801	0.402	ng	-0.02
31) Terphenyl-d14	19.447	244	10934	0.657	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	15353	2.671	ng	99

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

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