

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034519.D
 Acq On : 10 Oct 2024 13:54
 Operator : RC/JU
 Sample : P4263-01
 Misc :
 ALS Vial : 95 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE115D1-20240930

Quant Time: Oct 10 15:09:06 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	4083	0.400	ng	-0.02
7) Naphthalene-d8	10.137	136	11213	0.400	ng	-0.01
13) Acenaphthene-d10	14.019	164	6506	0.400	ng	-0.03
19) Phenanthrene-d10	16.784	188	12362	0.400	ng	#-0.02
29) Chrysene-d12	21.008	240	7593	0.400	ng	0.00
35) Perylene-d12	23.116	264	10565	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1167	0.101	ng	-0.01
5) Phenol-d6	6.593	99	489	0.036	ng	0.00
8) Nitrobenzene-d5	8.536	82	2115	0.247	ng	0.00
11) 2-Methylnaphthalene-d10	11.735	152	5352	0.323	ng	-0.02
14) 2,4,6-Tribromophenol	15.542	330	628	0.213	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	7327	0.277	ng	-0.02
27) Fluoranthene-d10	18.830	212	12847	0.412	ng	-0.02
31) Terphenyl-d14	19.447	244	9406	0.620	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	3.025	88	20866	4.088	ng	Qvalue 99

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

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