

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100924\
 Data File : BN034511.D
 Acq On : 09 Oct 2024 19:14
 Operator : RC/JU
 Sample : P4231-06
 Misc :
 ALS Vial : 84 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW10A-MW01S-20240926

Quant Time: Oct 10 01:46:56 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	3583	0.400	ng	-0.02
7) Naphthalene-d8	10.127	136	9742	0.400	ng	#-0.02
13) Acenaphthene-d10	14.026	164	5474	0.400	ng	-0.02
19) Phenanthrene-d10	16.790	188	10229	0.400	ng	-0.01
29) Chrysene-d12	21.006	240	6965	0.400	ng	0.00
35) Perylene-d12	23.116	264	7621	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1166	0.115	ng	-0.01
5) Phenol-d6	6.585	99	509	0.043	ng	-0.01
8) Nitrobenzene-d5	8.525	82	1921	0.258	ng	-0.01
11) 2-Methylnaphthalene-d10	11.735	152	4410	0.307	ng	-0.02
14) 2,4,6-Tribromophenol	15.537	330	662	0.267	ng	-0.01
15) 2-Fluorobiphenyl	12.647	172	6264	0.281	ng	-0.01
27) Fluoranthene-d10	18.832	212	10502	0.407	ng	-0.01
31) Terphenyl-d14	19.450	244	7249	0.521	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	3.032	88	768	0.171	ng	# 33

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

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