

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
 Data File : BN034470.D
 Acq On : 08 Oct 2024 14:21
 Operator : RC/JU
 Sample : P4226-01
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT180D2-20240924

Quant Time: Oct 08 14:56:40 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	7.380	152	5295	0.400	ng	-0.01
7) Naphthalene-d8	10.137	136	14843	0.400	ng	-0.01
13) Acenaphthene-d10	14.026	164	8348	0.400	ng	-0.02
19) Phenanthrene-d10	16.790	188	14510	0.400	ng	#-0.01
29) Chrysene-d12	21.015	240	11622	0.400	ng	0.00
35) Perylene-d12	23.125	264	12734	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.025	112	2066	0.137	ng	0.00
5) Phenol-d6	6.593	99	878	0.050	ng	0.00
8) Nitrobenzene-d5	8.536	82	3176	0.280	ng	0.00
11) 2-Methylnaphthalene-d10	11.742	152	7540	0.344	ng	-0.01
14) 2,4,6-Tribromophenol	15.549	330	1228	0.325	ng	0.00
15) 2-Fluorobiphenyl	12.647	172	10137	0.299	ng	-0.01
27) Fluoranthene-d10	18.836	212	14962	0.409	ng	0.00
31) Terphenyl-d14	19.459	244	10082	0.434	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	26629	4.023	ng	98
6) bis(2-Chloroethyl)ether	6.831	93	489	0.032	ng	87

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

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