

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
 Data File : BN034488.D
 Acq On : 09 Oct 2024 03:44
 Operator : RC/JU
 Sample : P4226-13
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D2-20240925

Quant Time: Oct 09 04:07:37 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	4127	0.400	ng	-0.01
7) Naphthalene-d8	10.137	136	10962	0.400	ng	-0.01
13) Acenaphthene-d10	14.030	164	6631	0.400	ng	-0.02
19) Phenanthrene-d10	16.784	188	12298	0.400	ng	#-0.02
29) Chrysene-d12	21.008	240	9089	0.400	ng	0.00
35) Perylene-d12	23.119	264	12420	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.025	112	1829	0.156	ng	0.00
5) Phenol-d6	6.585	99	780	0.057	ng	-0.01
8) Nitrobenzene-d5	8.536	82	2233	0.266	ng	0.00
11) 2-Methylnaphthalene-d10	11.738	152	5256	0.325	ng	-0.01
14) 2,4,6-Tribromophenol	15.542	330	1092	0.363	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	9985	0.370	ng	-0.02
27) Fluoranthene-d10	18.829	212	14438	0.465	ng	-0.02
31) Terphenyl-d14	19.452	244	11250	0.620	ng	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.032	88	24433	4.735	ng	97

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

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