

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100824\
 Data File : BN034491.D
 Acq On : 09 Oct 2024 05:32
 Operator : RC/JU
 Sample : P4226-16
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW9-MW-01D1-20240925

Quant Time: Oct 09 06:57:32 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	4544	0.400	ng	-0.01
7) Naphthalene-d8	10.137	136	12061	0.400	ng	-0.01
13) Acenaphthene-d10	14.030	164	6840	0.400	ng	-0.02
19) Phenanthrene-d10	16.796	188	12024	0.400	ng	0.00
29) Chrysene-d12	21.008	240	8021	0.400	ng	0.00
35) Perylene-d12	23.119	264	9925	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.026	112	2053	0.159	ng	0.00
5) Phenol-d6	6.586	99	905	0.060	ng	-0.01
8) Nitrobenzene-d5	8.536	82	2722	0.295	ng	0.00
11) 2-Methylnaphthalene-d10	11.742	152	5994	0.337	ng	-0.01
14) 2,4,6-Tribromophenol	15.555	330	1095	0.353	ng	0.00
15) 2-Fluorobiphenyl	12.651	172	12988	0.467	ng	0.00
27) Fluoranthene-d10	18.834	212	13801	0.455	ng	-0.01
31) Terphenyl-d14	19.452	244	10513	0.656	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	3.032	88	2121	0.373	ng	# 56

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

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