

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034146.D
 Acq On : 22 Sep 2024 04:47
 Operator : JU/RC
 Sample : P4111-04
 Misc :
 ALS Vial : 111 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW7-SP303-20240918

Quant Time: Sep 23 00:03:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 18 16:09:28 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.423	152	5454	0.400	ng	-0.01
7) Naphthalene-d8	10.180	136	14644	0.400	ng	#-0.02
13) Acenaphthene-d10	14.069	164	8075	0.400	ng	-0.01
19) Phenanthrene-d10	16.828	188	15340	0.400	ng	#-0.01
29) Chrysene-d12	21.051	240	10264	0.400	ng	# 0.00
35) Perylene-d12	23.169	264	9401	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	1970	0.127	ng	0.00
5) Phenol-d6	6.622	99	1222	0.068	ng	-0.01
8) Nitrobenzene-d5	8.568	82	3312	0.296	ng	0.00
11) 2-Methylnaphthalene-d10	11.784	152	6606	0.305	ng	0.00
14) 2,4,6-Tribromophenol	15.574	330	766	0.209	ng	-0.01
15) 2-Fluorobiphenyl	12.690	172	10186	0.310	ng	-0.01
27) Fluoranthene-d10	18.874	212	13674	0.353	ng	-0.01
31) Terphenyl-d14	19.491	244	8903	0.434	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.076	88	505	0.074	ng	# 1

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

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