

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092624\
 Data File : BN034256.D
 Acq On : 26 Sep 2024 22:11
 Operator : RC/JU
 Sample : P4164-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT180D1-20240923

Quant Time: Sep 27 01:03:06 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	7.401	152	5297	0.400	ng	0.00
7) Naphthalene-d8	10.159	136	14815	0.400	ng	0.01
13) Acenaphthene-d10	14.040	164	7830	0.400	ng	0.00
19) Phenanthrene-d10	16.796	188	14821	0.400	ng	# 0.00
29) Chrysene-d12	21.017	240	9879	0.400	ng	# 0.00
35) Perylene-d12	23.130	264	9250	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.040	112	1649	0.110	ng	0.00
5) Phenol-d6	6.607	99	644	0.037	ng	0.00
8) Nitrobenzene-d5	8.546	82	2771	0.245	ng	0.01
11) 2-Methylnaphthalene-d10	11.750	152	7475	0.342	ng	0.00
14) 2,4,6-Tribromophenol	15.555	330	642	0.181	ng	0.00
15) 2-Fluorobiphenyl	12.661	172	10180	0.320	ng	0.00
27) Fluoranthene-d10	18.848	212	13159	0.352	ng	0.00
31) Terphenyl-d14	19.466	244	8350	0.423	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.039	88	24316	3.672	ng	# 86

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

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