

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092624\
 Data File : BN034252.D
 Acq On : 26 Sep 2024 19:46
 Operator : RC/JU
 Sample : P4164-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE139D2-20240920

Quant Time: Sep 27 01:01:56 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	7.402	152	5235	0.400	ng	0.00
7) Naphthalene-d8	10.159	136	13458	0.400	ng	# 0.01
13) Acenaphthene-d10	14.051	164	6674	0.400	ng	0.00
19) Phenanthrene-d10	16.808	188	12156	0.400	ng	# 0.00
29) Chrysene-d12	21.026	240	8279	0.400	ng	0.01
35) Perylene-d12	23.133	264	8196	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.040	112	2401	0.162	ng	0.00
5) Phenol-d6	6.600	99	1033	0.060	ng	0.00
8) Nitrobenzene-d5	8.557	82	2600	0.253	ng	0.02
11) 2-Methylnaphthalene-d10	11.761	152	5821	0.293	ng	0.00
14) 2,4,6-Tribromophenol	15.555	330	245	0.081	ng	0.00
15) 2-Fluorobiphenyl	12.661	172	8353	0.308	ng	0.00
27) Fluoranthene-d10	18.848	212	10838	0.353	ng	0.00
31) Terphenyl-d14	19.466	244	6730	0.407	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.039	88	10278	1.570	ng	# 81

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

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