

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120123\
 Data File : BN028996.D
 Acq On : 02 Dec 2023 22:59
 Operator : MA/JU
 Sample : 05404-16
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE108D2-20231109

Quant Time: Dec 03 08:16:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN113023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 23:11:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.782	152	2306	0.400	ng	-0.01
7) Naphthalene-d8	10.573	136	6659	0.400	ng	#-0.01
13) Acenaphthene-d10	14.407	164	4309	0.400	ng	-0.01
19) Phenanthrene-d10	17.159	188	9765	0.400	ng	#-0.01
29) Chrysene-d12	21.356	240	7861	0.400	ng	#-0.02
35) Perylene-d12	23.678	264	7759	0.400	ng	#-0.03
System Monitoring Compounds						
4) 2-Fluorophenol	5.399	112	760	0.109	ng	-0.01
5) Phenol-d6	6.995	99	549	0.071	ng	-0.01
8) Nitrobenzene-d5	8.950	82	2056	0.311	ng	-0.02
11) 2-Methylnaphthalene-d10	12.159	152	3033	0.284	ng	-0.01
14) 2,4,6-Tribromophenol	15.905	330	750	0.258	ng	-0.01
15) 2-Fluorobiphenyl	13.038	172	5476	0.294	ng	-0.01
27) Fluoranthene-d10	19.190	212	9490	0.369	ng	-0.02
31) Terphenyl-d14	19.803	244	8786	0.395	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.333	88	4760	2.108	ng	# 91
6) bis(2-Chloroethyl)ether	7.240	93	347	0.053	ng	99

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

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