

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122220\
 Data File : BN013252.D
 Acq On : 22 Dec 2020 00:41
 Operator : CG/JU
 Sample : L5123-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW202D-20201213

Quant Time: Dec 22 01:35:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN121620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 18 16:11:20 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.338	152	624	0.40	ng	0.00
7) Naphthalene-d8	10.112	136	2303	0.40	ng	# 0.00
13) Acenaphthene-d10	13.991	164	1540	0.40	ng	-0.01
19) Phenanthrene-d10	16.763	188	3327	0.40	ng	#-0.01
29) Chrysene-d12	20.995	240	3822	0.40	ng	#-0.01
36) Perylene-d12	23.082	264	4454	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	4.962	112	282	0.11	ng	0.00
5) Phenol-d6	6.565	99	137	0.05	ng	0.02
8) Nitrobenzene-d5	8.541	82	602	0.25	ng	0.01
11) 2-Methylnaphthalene-d10	11.720	152	977	0.24	ng	0.00
14) 2,4,6-Tribromophenol	15.526	330	260	0.27	ng	-0.01
15) 2-Fluorobiphenyl	12.621	172	1649	0.26	ng	-0.01
27) Fluoranthene-d10	18.811	212	3984	0.36	ng	0.00
31) Terphenyl-d14	19.432	244	3499	0.36	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	2.925	88	752	0.69	ng	92
25) Phenanthrene	16.812	178	237	0.02	ng	# 82

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

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