

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122220\
 Data File : BN013247.D
 Acq On : 21 Dec 2020 21:41
 Operator : CG/JU
 Sample : L5123-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D-20201211

Quant Time: Dec 22 01:06:26 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	638	0.40	ng	0.00
7) Naphthalene-d8	10.112	136	2267	0.40	ng	# 0.00
13) Acenaphthene-d10	13.991	164	1494	0.40	ng	-0.01
19) Phenanthrene-d10	16.775	188	3126	0.40	ng	# 0.00
29) Chrysene-d12	21.003	240	3643	0.40	ng	# 0.00
36) Perylene-d12	23.086	264	4216	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.962	112	252	0.10	ng	0.00
5) Phenol-d6	6.557	99	136	0.05	ng	0.00
8) Nitrobenzene-d5	8.528	82	570	0.24	ng	0.00
11) 2-Methylnaphthalene-d10	11.720	152	923	0.23	ng	0.00
14) 2,4,6-Tribromophenol	15.526	330	235	0.25	ng	-0.01
15) 2-Fluorobiphenyl	12.608	172	1473	0.24	ng	-0.03
27) Fluoranthene-d10	18.813	212	3563	0.35	ng	0.00
31) Terphenyl-d14	19.429	244	3175	0.34	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	2.917	88	6157	5.53	ng	# 91
6) bis(2-Chloroethyl)ether	6.790	93	97	0.06	ng	# 83

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

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