

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122220\
 Data File : BN013267.D
 Acq On : 22 Dec 2020 11:35
 Operator : CG/JU
 Sample : L5123-21
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW179D-20201214

Quant Time: Dec 22 13:13:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN121620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 22 10:46:43 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.338	152	596	0.40	ng	0.00
7) Naphthalene-d8	10.112	136	2065	0.40	ng	# 0.01
13) Acenaphthene-d10	13.991	164	1391	0.40	ng	0.00
19) Phenanthrene-d10	16.775	188	3200	0.40	ng	# 0.01
29) Chrysene-d12	20.992	240	3605	0.40	ng	# 0.00
36) Perylene-d12	23.082	264	4123	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.962	112	287	0.12	ng	0.00
5) Phenol-d6	6.557	99	162	0.07	ng	0.02
8) Nitrobenzene-d5	8.540	82	465	0.22	ng	0.04
11) 2-Methylnaphthalene-d10	11.719	152	869	0.24	ng	0.01
14) 2,4,6-Tribromophenol	15.526	330	242	0.28	ng	0.00
15) 2-Fluorobiphenyl	12.621	172	1318	0.23	ng	0.01
27) Fluoranthene-d10	18.813	212	3682	0.35	ng	0.00
31) Terphenyl-d14	19.429	244	3417	0.37	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.933	88	150	0.14	ng	# 58
32) Benzo(a)anthracene	20.982	228	389	0.03	ng	# 67
33) Chrysene	21.035	228	425	0.03	ng	# 75
37) Benzo(b)fluoranthene	22.477	252	487	0.03	ng	# 1
38) Benzo(k)fluoranthene	22.514	252	425	0.03	ng	# 1
39) Benzo(a)pyrene	22.997	252	399	0.03	ng	# 1

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

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