

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120321\
 Data File : BN017705.D
 Acq On : 03 Dec 2021 23:56
 Operator : CG/JU
 Sample : M4917-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-16R

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021
 Supervised By :Sohil Jodhani 12/10/2021

Quant Time: Dec 04 01:01:24 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN120321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 03 23:07:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.754	152	25285	0.40	ng	0.00
7) Naphthalene-d8	10.535	136	105936	0.40	ng	# 0.00
13) Acenaphthene-d10	14.372	164	69761	0.40	ng	0.00
19) Phenanthrene-d10	17.116	188	142365	0.40	ng	0.00
29) Chrysene-d12	21.303	240	147356	0.40	ng	# 0.00
35) Perylene-d12	23.609	264	141568	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.366	112	6592	0.11	ng	0.00
5) Phenol-d6	6.934	99	3404	0.06	ng	0.00
8) Nitrobenzene-d5	8.900	82	15466m	0.30	ng	0.00
11) 2-Methylnaphthalene-d10	12.124	152	38871	0.21	ng	0.00
14) 2,4,6-Tribromophenol	15.856	330	10917	0.38	ng	-0.01
15) 2-Fluorobiphenyl	13.001	172	69105	0.26	ng	0.00
27) Fluoranthene-d10	19.169	212	113062	0.24	ng	0.02
31) Terphenyl-d14	19.749	244	115986	0.35	ng	0.00
Target Compounds						
9) Naphthalene	10.585	128	5434	0.02	ng	# 89
24) Pentachlorophenol	16.776	266	930	0.30	ng	93
34) Bis(2-ethylhexyl)phtha...	21.229	149	30922	0.22	ng	# 96

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

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