

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091422\
 Data File : BN021751.D
 Acq On : 14 Sep 2022 17:32
 Operator : CG/JU
 Sample : N4549-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-46

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/15/2022
 Supervised By : Jagrut Upadhyay 09/15/2022

Quant Time: Sep 15 02:11:22 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 14 01:55:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.054	152	4819	0.400	ng	0.00
7) Naphthalene-d8	10.887	136	16397	0.400	ng	0.00
13) Acenaphthene-d10	14.707	164	9922	0.400	ng	0.00
19) Phenanthrene-d10	17.456	188	21880	0.400	ng	0.00
29) Chrysene-d12	21.647	240	17381	0.400	ng	# 0.00
35) Perylene-d12	24.160	264	14317	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.584	112	2298	0.182	ng	0.00
5) Phenol-d6	7.209	99	1967	0.123	ng	0.00
8) Nitrobenzene-d5	9.233	82	4535m	0.335	ng	0.00
11) 2-Methylnaphthalene-d10	12.471	152	16128	0.390	ng	0.00
14) 2,4,6-Tribromophenol	16.197	330	1918	0.423	ng	-0.01
15) 2-Fluorobiphenyl	13.339	172	18996	0.469	ng	0.00
27) Fluoranthene-d10	19.494	212	25983	0.441	ng	0.00
31) Terphenyl-d14	20.080	244	20394	0.513	ng	0.00
Target Compounds						Qvalue
34) Bis(2-ethylhexyl)phtha...	21.549	149	3415	0.077	ng	# 100

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

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