

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091422\
 Data File : BN021748.D
 Acq On : 14 Sep 2022 15:40
 Operator : CG/JU
 Sample : N4549-16
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-60

Manual Integrations
 APPROVED

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

Quant Time: Sep 15 02:10:30 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	8.055	152	5337	0.400	ng	0.00
7) Naphthalene-d8	10.888	136	18162	0.400	ng	0.00
13) Acenaphthene-d10	14.707	164	11219	0.400	ng	0.00
19) Phenanthrene-d10	17.456	188	23721	0.400	ng	0.00
29) Chrysene-d12	21.652	240	17507	0.400	ng	0.00
35) Perylene-d12	24.158	264	13718	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.585	112	1926	0.138	ng	0.00
5) Phenol-d6	7.210	99	1527	0.086	ng	0.00
8) Nitrobenzene-d5	9.233	82	5129m	0.342	ng	0.00
11) 2-Methylnaphthalene-d10	12.471	152	14918	0.326	ng	0.00
14) 2,4,6-Tribromophenol	16.197	330	1993	0.389	ng	-0.01
15) 2-Fluorobiphenyl	13.339	172	17840	0.389	ng	0.00
27) Fluoranthene-d10	19.490	212	25850	0.405	ng	0.00
31) Terphenyl-d14	20.076	244	17050	0.426	ng	0.00
Target Compounds						
34) Bis(2-ethylhexyl)phtha...	21.544	149	14099	0.314	ng	Qvalue 99

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

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