

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091422\
 Data File : BN021763.D
 Acq On : 15 Sep 2022 01:37
 Operator : CG/JU
 Sample : N4549-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-48

Manual Integrations
 APPROVED

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

Quant Time: Sep 15 02:15:07 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	5138	0.400	ng	0.00
7) Naphthalene-d8	10.888	136	18005	0.400	ng	# 0.00
13) Acenaphthene-d10	14.707	164	16477	0.400	ng	# 0.00
19) Phenanthrene-d10	17.456	188	23353	0.400	ng	# 0.00
29) Chrysene-d12	21.647	240	16585	0.400	ng	# 0.00
35) Perylene-d12	24.160	264	12636	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.584	112	2704	0.201	ng	0.00
5) Phenol-d6	7.209	99	2208	0.130	ng	0.00
8) Nitrobenzene-d5	9.222	82	5247	0.352	ng	-0.01
11) 2-Methylnaphthalene-d10	12.471	152	9989	0.220	ng	0.00
14) 2,4,6-Tribromophenol	16.197	330	2592	0.344	ng	-0.01
15) 2-Fluorobiphenyl	13.339	172	19641	0.292	ng	0.00
27) Fluoranthene-d10	19.490	212	24717	0.393	ng	0.00
31) Terphenyl-d14	20.080	244	21594	0.570	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.930	128	4893	0.092	ng	# 35
16) Acenaphthylene	14.429	152	9204	0.117	ng	# 1
17) Acenaphthene	14.771	154	40549	0.754	ng	96
18) Fluorene	15.747	166	6620	0.093	ng	# 91
34) Bis(2-ethylhexyl)phtha...	21.540	149	6263	0.147	ng	# 86

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

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