

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091422\
 Data File : BN021764.D
 Acq On : 15 Sep 2022 02:13
 Operator : CG/JU
 Sample : N4549-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-47

Quant Time: Sep 15 02:57:50 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	5251	0.400	ng	0.00
7) Naphthalene-d8	10.888	136	18156	0.400	ng	# 0.00
13) Acenaphthene-d10	14.707	164	18413	0.400	ng	# 0.00
19) Phenanthrene-d10	17.456	188	21503	0.400	ng	# 0.00
29) Chrysene-d12	21.652	240	14574	0.400	ng	# 0.00
35) Perylene-d12	24.158	264	10579	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.584	112	4087	0.298	ng	0.00
5) Phenol-d6	7.209	99	2199	0.126	ng	0.00
8) Nitrobenzene-d5	9.222	82	5014	0.334	ng	-0.01
11) 2-Methylnaphthalene-d10	12.471	152	13638	0.298	ng	0.00
14) 2,4,6-Tribromophenol	16.197	330	2354	0.280	ng	-0.01
15) 2-Fluorobiphenyl	13.339	172	19567	0.260	ng	0.00
27) Fluoranthene-d10	19.494	212	22997	0.397	ng	0.00
31) Terphenyl-d14	20.084	244	18343	0.551	ng	0.00
Target Compounds						
9) Naphthalene	10.930	128	3731	0.070	ng	# 13
17) Acenaphthene	14.771	154	11761	0.196	ng	92
34) Bis(2-ethylhexyl)phtha...	21.544	149	3775	0.101	ng	# 41

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

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