

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091721\
 Data File : BN016324.D
 Acq On : 18 Sep 2021 08:32
 Operator : CG/JU
 Sample : M3791-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SW-RR-02-(1.0)

Manual Integrations
 APPROVED

mohammad
 9/20/2021 4:22:30 PM

Quant Time: Sep 20 00:55:49 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 17 23:59:14 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.689	152	41508	0.40	ng	0.00
7) Naphthalene-d8	10.457	136	185811	0.40	ng	0.00
13) Acenaphthene-d10	14.307	164	110988	0.40	ng	0.00
19) Phenanthrene-d10	17.054	188	221695	0.40	ng	#-0.01
29) Chrysene-d12	21.259	240	239161	0.40	ng	# 0.00
35) Perylene-d12	23.526	264	255662	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.319	112	7595	0.07	ng	0.02
5) Phenol-d6	6.868	99	5241	0.04	ng	0.00
8) Nitrobenzene-d5	8.822	82	21807	0.20	ng	-0.01
11) 2-Methylnaphthalene-d10	12.047	152	49619	0.17	ng	0.00
14) 2,4,6-Tribromophenol	15.804	330	8721	0.27	ng	0.00
15) 2-Fluorobiphenyl	12.937	172	68787	0.16	ng	0.00
27) Fluoranthene-d10	19.098	212	121179	0.19	ng	0.00
31) Terphenyl-d14	19.708	244	105637	0.19	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.508	128	1122958	2.24	ng	100
12) 2-Methylnaphthalene	12.123	142	24997	0.08	ng	98
34) Bis(2-ethylhexyl)phtha...	21.196	149	47292	0.11	ng	99

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

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