

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
 Data File : BM026854.D
 Acq On : 23 Jul 2020 16:54
 Operator : JU/CG
 Sample : L3381-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 357-GRID-A4-A6

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:16:26 PM

Quant Time: Jul 24 04:03:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 22 14:16:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.28	152	96133	20.00	ng	-0.02
21) Naphthalene-d8	10.04	136	373664	20.00	ng	-0.01
39) Acenaphthene-d10	13.95	164	216296	20.00	ng	-0.01
64) Phenanthrene-d10	16.71	188	447240	20.00	ng	-0.01
76) Chrysene-d12	20.94	240	483024	20.00	ng	0.00
86) Perylene-d12	23.01	264	511978	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.92	112	474891	82.80	ng	-0.02
7) Phenol-d6	6.50	99	640081	70.04	ng	-0.02
23) Nitrobenzene-d5	8.43	82	433039	49.44	ng	-0.02
42) 2,4,6-Tribromophenol	15.46	330	218794	69.44	ng	-0.01
45) 2-Fluorobiphenyl	12.55	172	829071	50.97	ng	-0.01
79) Terphenyl-d14	19.38	244	1169621	47.65	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.42	163	175307	9.476	ng	97
78) Pyrene	19.15	202	68442	2.249	ng	97
84) Bis(2-ethylhexyl)phthalate	20.89	149	211804	9.683	ng	98
88) Benzo(b)fluoranthene	22.41	252	68834m	2.040	ng	

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

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