

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
 Data File : BM026858.D
 Acq On : 23 Jul 2020 19:21
 Operator : JU/CG
 Sample : L3381-16
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 365-GRID-B6-B8

Quant Time: Jul 24 01:59:19 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	104108	20.00	ng	-0.02
21) Naphthalene-d8	10.03	136	385436	20.00	ng	-0.02
39) Acenaphthene-d10	13.95	164	193005	20.00	ng	-0.01
64) Phenanthrene-d10	16.71	188	356940	20.00	ng	-0.01
76) Chrysene-d12	20.94	240	378875	20.00	ng	0.00
86) Perylene-d12	23.01	264	413222	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.93	112	540600	87.03	ng	-0.02
7) Phenol-d6	6.50	99	714907	72.24	ng	-0.02
23) Nitrobenzene-d5	8.43	82	469530	51.97	ng	-0.02
42) 2,4,6-Tribromophenol	15.46	330	181751	64.65	ng	-0.01
45) 2-Fluorobiphenyl	12.55	172	825600	56.88	ng	-0.02
79) Terphenyl-d14	19.38	244	913412	47.45	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.42	163	170788	10.345	ng	99
71) Phenanthrene	16.75	178	54394	2.581	ng	99
75) Fluoranthene	18.79	202	137303	5.290	ng	98
78) Pyrene	19.15	202	144955	6.073	ng	99
81) Benzo(a)anthracene	20.92	228	77109	3.030	ng	97
83) Chrysene	20.98	228	82881	3.346	ng	97
88) Benzo(b)fluoranthene	22.41	252	110650	4.064	ng	# 87
90) Benzo(a)pyrene	22.92	252	76490	3.026	ng	# 87
92) Benzo(g,h,i)perylene	25.60	276	58035	2.536	ng	# 93

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

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