

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060319\
 Data File : BM020686.D
 Acq On : 04 Jun 2019 00:18
 Operator : HP/JU
 Sample : K3139-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-1A-S

Manual Integrations
 APPROVED

mohammad
 6/5/2019 8:46:09 AM

Quant Time: Jun 04 07:04:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 31 15:48:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	224375	20.00	ng	0.00
21) Naphthalene-d8	10.62	136	838251	20.00	ng	-0.01
39) Acenaphthene-d10	14.46	164	414624	20.00	ng	-0.01
64) Phenanthrene-d10	17.20	188	858677	20.00	ng	-0.01
76) Chrysene-d12	21.38	240	893559	20.00	ng	-0.02
87) Perylene-d12	23.66	264	1080114	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1676436	131.43	ng	0.00
7) Phenol-d6	6.99	99	2273757	121.06	ng	0.00
23) Nitrobenzene-d5	8.99	82	1370668	84.72	ng	-0.01
42) 2,4,6-Tribromophenol	15.95	330	642055	124.98	ng	-0.01
45) 2-Fluorobiphenyl	13.08	172	2783651	89.25	ng	-0.01
79) Terphenyl-d14	19.83	244	3753358	70.26	ng	-0.02
Target Compounds						
50) Dimethylphthalate	13.92	163	162533	4.588	ng	99
71) Phenanthrene	17.24	178	533952	10.970	ng	98
75) Fluoranthene	19.26	202	502321	9.280	ng	97
78) Pyrene	19.62	202	389279	5.677	ng	99
81) Benzo(a)anthracene	21.36	228	264224	4.246	ng	99
83) Chrysene	21.42	228	225937	3.820	ng	97
88) Benzo(b)fluoranthene	22.98	252	262803m	3.681	ng	
90) Benzo(a)pyrene	23.56	252	162408	2.508	ng	99

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

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