

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM083019\
 Data File : BM022441.D
 Acq On : 30 Aug 2019 14:35
 Operator : HP/JU
 Sample : K4574-14
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 S702B-N-1.0-1.5-082719-FD

Quant Time: Aug 30 15:45:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 16:51:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	16678	20.00	ng	-0.01
21) Naphthalene-d8	10.30	136	53109	20.00	ng	0.00
39) Acenaphthene-d10	14.18	164	32134	20.00	ng	0.00
64) Phenanthrene-d10	16.95	188	72475	20.00	ng	0.00
76) Chrysene-d12	21.17	240	99489	20.00	ng	0.00
87) Perylene-d12	23.36	264	131282	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.14	112	122441	128.87	ng	0.00
7) Phenol-d6	6.72	99	152014	119.64	ng	0.00
23) Nitrobenzene-d5	8.69	82	105746	87.21	ng	0.00
42) 2,4,6-Tribromophenol	15.70	330	114934	174.79	ng	0.00
45) 2-Fluorobiphenyl	12.79	172	251903	88.81	ng	0.00
79) Terphenyl-d14	19.60	244	865282	114.78	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.66	163	16035	5.737	ng	98
75) Fluoranthene	19.03	202	22406	3.983	ng	96
78) Pyrene	19.39	202	20517	3.026	ng	96
81) Benzo(a)anthracene	21.15	228	16421	2.300	ng	96
83) Chrysene	21.21	228	15508	2.282	ng	97
88) Benzo(b)fluoranthene	22.71	252	25125	2.755	ng	# 84
90) Benzo(a)pyrene	23.26	252	18235	2.253	ng	# 90

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

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