

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM020320\
 Data File : BM024698.D
 Acq On : 03 Feb 2020 14:34
 Operator : CG/JU
 Sample : L1343-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 NW-SP

Manual Integrations
 APPROVED

mohammad
 2/4/2020 6:12:18 PM

Quant Time: Feb 04 03:26:53 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 23 14:57:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	208992	20.00	ng	-0.02
21) Naphthalene-d8	10.18	136	756579	20.00	ng	-0.02
39) Acenaphthene-d10	14.07	164	508488	20.00	ng	-0.02
64) Phenanthrene-d10	16.83	188	1173070	20.00	ng	-0.02
76) Chrysene-d12	21.03	240	1232780	20.00	ng	-0.02
87) Perylene-d12	23.14	264	1451750	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.07	112	1346694	122.54	ng	-0.01
7) Phenol-d6	6.62	99	1684456	111.27	ng	-0.02
23) Nitrobenzene-d5	8.56	82	1089165	70.60	ng	-0.02
42) 2,4,6-Tribromophenol	15.57	330	993993	120.17	ng	-0.02
45) 2-Fluorobiphenyl	12.69	172	2948824	78.84	ng	-0.02
79) Terphenyl-d14	19.49	244	5402527	81.82	ng	-0.02
Target Compounds						
50) Dimethylphthalate	13.54	163	84265	2.049	ng	Qvalue 100
71) Phenanthrene	16.87	178	220346	3.469	ng	98
75) Fluoranthene	18.89	202	623385	7.863	ng	98
78) Pyrene	19.26	202	464479	5.715	ng	99
83) Chrysene	21.07	228	181811	2.227	ng	98
88) Benzo(b)fluoranthene	22.53	252	258478m	2.745	ng	

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

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