

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
 Data File : BF118389.D
 Acq On : 30 Dec 2019 19:05
 Operator : JU
 Sample : K6442-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K615-WC-1

Manual Integrations
 APPROVED

mohammad
 12/31/2019 12:04:17 PM

Quant Time: Dec 31 01:36:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	125640	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	488819	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	266454	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	489724	20.00	ng	0.00
76) Chrysene-d12	14.05	240	304926	20.00	ng	0.00
87) Perylene-d12	15.54	264	328807	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	689752	93.71	ng	0.00
7) Phenol-d6	6.49	99	875268	95.86	ng	0.00
23) Nitrobenzene-d5	7.42	82	526026	61.55	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	313030	95.13	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1146306	65.67	ng	0.00
79) Terphenyl-d14	12.99	244	1297608	70.60	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.62	163	90424	4.521	ng	99
71) Phenanthrene	11.42	178	74612	2.795	ng	98
75) Fluoranthene	12.62	202	101211	3.756	ng	99
78) Pyrene	12.84	202	89180	3.581	ng	99
81) Benzo(a)anthracene	14.04	228	44642	2.092	ng	98
88) Benzo(b)fluoranthene	15.10	252	48803m	2.236	ng	
90) Benzo(a)pyrene	15.48	252	40799	2.122	ng	94
92) Benzo(g,h,i)perylene	17.51	276	35335	2.192	ng #	94

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

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