

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022119\
 Data File : BF112648.D
 Acq On : 21 Feb 2019 13:14
 Operator : JU/SJ
 Sample : K1546-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR-200034

Manual Integrations
 APPROVED

Sohil
 2/21/2019 3:59:49 PM

Quant Time: Feb 21 15:37:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 21 12:57:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	86694	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	340335	20.00	ng	0.00
39) Acenaphthene-d10	9.86	164	178101	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	363371	20.00	ng	0.00
76) Chrysene-d12	14.00	240	287053	20.00	ng	0.00
87) Perylene-d12	15.47	264	225074	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	584804	143.28	ng	0.00
7) Phenol-d6	6.47	99	726920	128.17	ng	0.00
23) Nitrobenzene-d5	7.39	82	477397	80.82	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	269549	130.03	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	955541	75.88	ng	0.00
79) Terphenyl-d14	12.93	244	1157342	75.94	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.57	163	69941	5.054	ng	97
75) Fluoranthene	12.57	202	78778	3.888	ng	96
78) Pyrene	12.80	202	72267	3.636	ng	98
83) Chrysene	14.02	228	35258m	2.189	ng	
88) Benzo(b)fluoranthene	15.04	252	38094m	2.830	ng	

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

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