

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
 Data File : BF112277.D
 Acq On : 28 Jan 2019 16:25
 Operator : JU/SJ
 Sample : K1261-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-C05

Manual Integrations
 APPROVED

mohammad
 1/29/2019 9:35:11 AM

Quant Time: Jan 28 16:59:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	229549	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	902845	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	451520	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	900546	20.00	ng	0.00
76) Chrysene-d12	14.00	240	693238	20.00	ng	0.00
87) Perylene-d12	15.47	264	532421	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1662232	153.81	ng	0.00
7) Phenol-d6	6.49	99	2135005	142.17	ng	0.00
23) Nitrobenzene-d5	7.40	82	1384330	88.35	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	714747	136.00	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	2539418	79.54	ng	0.00
79) Terphenyl-d14	12.95	244	2711750	73.68	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	184424	5.257	ng	99
75) Fluoranthene	12.57	202	205583	4.094	ng	96
78) Pyrene	12.80	202	187107	3.898	ng	97
81) Benzo(a)anthracene	13.99	228	88124	2.162	ng	99
88) Benzo(b)fluoranthene	15.04	252	73584m	2.311	ng	

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

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