

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
 Data File : BF113554.D
 Acq On : 3 Apr 2019 20:16
 Operator : JU/SJ
 Sample : K2213-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S9A-S9B-S9C

Manual Integrations
 APPROVED

Sohil
 4/4/2019 10:57:49 AM

Quant Time: Apr 04 05:33:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	121711	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	422673	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	175853	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	326929	20.00	ng	0.00
76) Chrysene-d12	13.84	240	317124	20.00	ng	0.00
87) Perylene-d12	15.25	264	279469	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	784446	109.90	ng	0.00
7) Phenol-d6	6.33	99	986383	112.13	ng	0.00
23) Nitrobenzene-d5	7.25	82	582116	80.67	ng	0.00
42) 2,4,6-Tribromophenol	10.51	330	232867	110.24	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	969318	88.74	ng	0.00
79) Terphenyl-d14	12.79	244	943115	60.55	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.44	163	65520	5.054	ng	99
75) Fluoranthene	12.42	202	68973	3.962	ng	98
78) Pyrene	12.64	202	71168	2.949	ng	99
81) Benzo(a)anthracene	13.83	228	48863	2.671	ng	99
83) Chrysene	13.86	228	56627	3.054	ng	98
88) Benzo(b)fluoranthene	14.85	252	65346m	3.820	ng	

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

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