

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
 Data File : BF113553.D
 Acq On : 3 Apr 2019 19:47
 Operator : JU/SJ
 Sample : K2213-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S9C(2-10)COMP

Manual Integrations
 APPROVED

Sohil
 4/4/2019 10:57:47 AM

Quant Time: Apr 04 05:28:27 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	137887	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	472609	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	199039	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	354707	20.00	ng	0.00
76) Chrysene-d12	13.84	240	330557	20.00	ng	0.00
87) Perylene-d12	15.25	264	288629	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1014946	125.52	ng	0.00
7) Phenol-d6	6.34	99	1293390	129.78	ng	0.00
23) Nitrobenzene-d5	7.25	82	775282	96.09	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	307202	128.49	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	1348265	109.05	ng	0.00
79) Terphenyl-d14	12.79	244	1293966	79.70	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.44	163	93466	6.370	ng	99
75) Fluoranthene	12.42	202	142777	7.559	ng	98
78) Pyrene	12.64	202	159178	6.329	ng	98
81) Benzo(a)anthracene	13.83	228	80028	4.197	ng	100
83) Chrysene	13.86	228	89381	4.624	ng	99
88) Benzo(b)fluoranthene	14.85	252	113635m	6.432	ng	
89) Benzo(k)fluoranthene	14.87	252	34223m	2.158	ng	
90) Benzo(a)pyrene	15.19	252	42936	2.735	ng	# 93

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

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