

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
 Data File : BF113552.D
 Acq On : 3 Apr 2019 19:18
 Operator : JU/SJ
 Sample : K2213-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S9A(2-10)COMP

Manual Integrations
 APPROVED

Sohil
 4/4/2019 5:39:23 PM

Quant Time: Apr 04 05:09:44 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	512958	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	240260	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	449566	20.00	ng	0.00
76) Chrysene-d12	13.84	240	361490	20.00	ng	0.00
87) Perylene-d12	15.25	264	305897	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	840285	103.92	ng	0.00
7) Phenol-d6	6.33	99	1094734	109.84	ng	0.00
23) Nitrobenzene-d5	7.25	82	646035	73.77	ng	0.00
42) 2,4,6-Tribromophenol	10.51	330	297530	103.09	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	1212129	81.22	ng	0.00
79) Terphenyl-d14	12.79	244	1200739	67.63	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.44	163	83528	4.716	ng	100
72) Anthracene	11.28	178	47904	2.108	ng	97
75) Fluoranthene	12.42	202	229441	9.585	ng	98
78) Pyrene	12.64	202	215566	7.837	ng	98
81) Benzo(a)anthracene	13.83	228	94804	4.546	ng	100
83) Chrysene	13.86	228	126139	5.967	ng	99
88) Benzo(b)fluoranthene	14.85	252	112917m	6.030	ng	
89) Benzo(k)fluoranthene	14.87	252	34048m	2.026	ng	
90) Benzo(a)pyrene	15.19	252	35391	2.127	ng	# 87

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

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