

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
 Data File : BF113899.D
 Acq On : 23 Apr 2019 00:55
 Operator : JU/SJ
 Sample : K2201-05 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-041819-C

Manual Integrations
 APPROVED

Sohil
 4/23/2019 4:08:17 PM

Quant Time: Apr 23 06:47:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	164421	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	577797	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	284056	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	609114	20.00	ng	0.00
76) Chrysene-d12	14.06	240	575565	20.00	ng	-0.01
87) Perylene-d12	15.54	264	460150	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	428043	44.53	ng	0.00
7) Phenol-d6	6.52	99	523162	43.38	ng	0.00
23) Nitrobenzene-d5	7.45	82	287982	30.77	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	169213	48.90	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	603020	33.20	ng	0.00
79) Terphenyl-d14	13.00	244	898785	32.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	49193	2.395	ng	Qvalue 100
71) Phenanthrene	11.44	178	126253	4.124	ng	98
75) Fluoranthene	12.63	202	323461	9.685	ng	99
78) Pyrene	12.86	202	317012	7.876	ng	98
81) Benzo(a)anthracene	14.04	228	215538m	6.356	ng	
83) Chrysene	14.08	228	178485	5.404	ng	97
86) Indeno(1,2,3-cd)pyrene	17.02	276	88244	2.821	ng	# 94
88) Benzo(b)fluoranthene	15.10	252	218677m	7.511	ng	
89) Benzo(k)fluoranthene	15.13	252	73165m	2.922	ng	
90) Benzo(a)pyrene	15.47	252	148356	5.891	ng	97
92) Benzo(g,h,i)perylene	17.47	276	80120	3.359	ng	95

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

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