

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021319\
 Data File : BF112533.D
 Acq On : 13 Feb 2019 12:32
 Operator : JU/SJ
 Sample : K1343-13 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-021119-G

Manual Integrations
 APPROVED

Sohil
 2/14/2019 11:36:07 AM

Quant Time: Feb 14 06:16:47 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	6.83	152	93515	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	350376	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	152440	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	256769	20.00	ng	0.00
76) Chrysene-d12	14.01	240	225309	20.00	ng	0.00
87) Perylene-d12	15.49	264	232593	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	286735	65.13	ng	-0.01
7) Phenol-d6	6.48	99	358482	58.60	ng	-0.01
23) Nitrobenzene-d5	7.40	82	221837	36.48	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	86373	48.68	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	420742	39.04	ng	-0.02
79) Terphenyl-d14	12.94	244	364415	30.46	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	34653	2.926	ng	Qvalue # 98
71) Phenanthrene	11.39	178	66871	5.009	ng	98
75) Fluoranthene	12.58	202	93476	6.529	ng	98
78) Pyrene	12.82	202	90123	5.777	ng	95
81) Benzo(a)anthracene	14.00	228	43684	3.298	ng	97
83) Chrysene	14.04	228	42724m	3.379	ng	
88) Benzo(b)fluoranthene	15.06	252	55159m	3.965	ng	
90) Benzo(a)pyrene	15.43	252	38303	3.060	ng	# 85
92) Benzo(g,h,i)perylene	17.46	276	19638	2.063	ng	# 83

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

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