

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012419\
 Data File : BF112216.D
 Acq On : 24 Jan 2019 18:38
 Operator : JU/SJ
 Sample : K1223-04 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-P

Manual Integrations
 APPROVED

Sohil
 1/25/2019 10:05:16 AM

Quant Time: Jan 25 00:34:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 19 01:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	167099	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	801898	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	354487	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	578964	20.00	ng	0.00
76) Chrysene-d12	14.03	240	441549	20.00	ng	0.00
87) Perylene-d12	15.50	264	310875	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	585162	63.56	ng	0.00
7) Phenol-d6	6.50	99	636577	55.27	ng	0.00
23) Nitrobenzene-d5	7.42	82	438297	37.76	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	184502	48.25	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	794709	32.87	ng	0.00
79) Terphenyl-d14	12.96	244	607609	29.27	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	66597	2.576	ng	Qvalue # 98
71) Phenanthrene	11.40	178	320457	10.998	ng	97
72) Anthracene	11.46	178	81060	2.746	ng	98
75) Fluoranthene	12.60	202	376625	12.243	ng	97
78) Pyrene	12.83	202	326096	11.194	ng	98
81) Benzo(a)anthracene	14.02	228	152012	6.004	ng	97
83) Chrysene	14.05	228	128975	5.384	ng	97
86) Indeno(1,2,3-cd)pyrene	17.00	276	46811	2.118	ng	# 93
88) Benzo(b)fluoranthene	15.07	252	136071m	7.711	ng	
89) Benzo(k)fluoranthene	15.10	252	42915m	2.513	ng	
90) Benzo(a)pyrene	15.44	252	85922	5.309	ng	# 85
92) Benzo(a,h,i)perylene	17.45	276	36163	2.554	ng	# 90

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

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