

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043019\
 Data File : BF114044.D
 Acq On : 30 Apr 2019 16:10
 Operator : JU/SJ
 Sample : K2194-03 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-042919-B

Manual Integrations
 APPROVED

Sohil
 5/1/2019 11:58:47 AM

Quant Time: May 01 05:20:48 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	6.89	152	124782	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	466894	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	223280	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	383943	20.00	ng	0.00
76) Chrysene-d12	14.05	240	305594	20.00	ng	-0.02
87) Perylene-d12	15.54	264	233573	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	392410	53.79	ng	0.00
7) Phenol-d6	6.52	99	521551	56.98	ng	0.00
23) Nitrobenzene-d5	7.45	82	294365	38.93	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	138265	50.83	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	606748	42.50	ng	-0.01
79) Terphenyl-d14	13.00	244	610385	40.95	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.63	163	84787	5.251	ng	98
71) Phenanthrene	11.43	178	137399	7.121	ng	98
75) Fluoranthene	12.63	202	232711	11.054	ng	97
78) Pyrene	12.86	202	212016	9.921	ng	98
81) Benzo(a)anthracene	14.04	228	120438	6.689	ng	95
83) Chrysene	14.07	228	112922	6.440	ng	96
86) Indeno(1,2,3-cd)pyrene	17.03	276	38792	2.335	ng	93
88) Benzo(b)fluoranthene	15.10	252	110280m	7.462	ng	
89) Benzo(k)fluoranthene	15.13	252	32592m	2.564	ng	
90) Benzo(a)pyrene	15.47	252	69601	5.445	ng	# 86
92) Benzo(a,h,i)perylene	17.47	276	39367	3.251	ng	# 86

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

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