

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042319\
 Data File : BF113938.D
 Acq On : 23 Apr 2019 23:12
 Operator : JU/SJ
 Sample : K2204-03 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-042219

Manual Integrations
 APPROVED

Sohil
 4/25/2019 6:02:26 PM

Quant Time: Apr 25 08:13:54 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	125284	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	428339	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	185691	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	353640	20.00	ng	0.00
76) Chrysene-d12	14.05	240	382614	20.00	ng	-0.02
87) Perylene-d12	15.53	264	310302	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	451961	61.70	ng	0.00
7) Phenol-d6	6.52	99	570198	62.04	ng	0.00
23) Nitrobenzene-d5	7.45	82	320102	46.14	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	134271	59.36	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	609020	51.30	ng	0.00
79) Terphenyl-d14	13.00	244	724808	38.84	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	77317	5.757	ng	100
71) Phenanthrene	11.44	178	93289	5.249	ng	96
75) Fluoranthene	12.63	202	208673	10.762	ng	98
78) Pyrene	12.86	202	205656	7.686	ng	97
81) Benzo(a)anthracene	14.04	228	113253	5.024	ng	97
83) Chrysene	14.08	228	116040	5.285	ng	97
86) Indeno(1,2,3-cd)pyrene	17.02	276	41807	2.010	ng	96
88) Benzo(b)fluoranthene	15.10	252	128856m	6.563	ng	
89) Benzo(k)fluoranthene	15.13	252	33794m	2.001	ng	
90) Benzo(a)pyrene	15.47	252	77354	4.555	ng	98
92) Benzo(g,h,i)perylene	17.47	276	38675	2.404	ng	# 93

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

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