

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060519\
 Data File : BF114843.D
 Acq On : 6 Jun 2019 2:36
 Operator : HP/JU
 Sample : K3151-05 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-060419-C

Manual Integrations
 APPROVED

mohammad
 6/6/2019 4:50:48 PM

Quant Time: Jun 06 07:59:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 04 16:57:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	348192	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	1337613	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	649275	20.00	ng	0.00
64) Phenanthrene-d10	11.17	188	1168758	20.00	ng	0.00
76) Chrysene-d12	13.80	240	766179	20.00	ng	0.00
87) Perylene-d12	15.18	264	633310	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	457937	22.98	ng	0.00
7) Phenol-d6	6.30	99	782203	32.26	ng	0.00
23) Nitrobenzene-d5	7.22	82	442475	20.36	ng	-0.01
42) 2,4,6-Tribromophenol	10.48	330	257152	36.75	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	1111791	32.08	ng	0.00
79) Terphenyl-d14	12.76	244	1204851	30.58	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.42	163	179286	3.781	ng	100
71) Phenanthrene	11.20	178	877936	15.735	ng	96
72) Anthracene	11.24	178	247562	4.216	ng	94
75) Fluoranthene	12.38	202	1615792	27.824	ng	94
78) Pyrene	12.61	202	1652441	24.861	ng	97
81) Benzo(a)anthracene	13.79	228	826672	14.001	ng	99
83) Chrysene	13.82	228	716484	12.475	ng	96
86) Indeno(1,2,3-cd)pyrene	16.49	276	287014	4.380	ng	96
88) Benzo(b)fluoranthene	14.80	252	699969m	17.361	ng	
89) Benzo(k)fluoranthene	14.82	252	256872m	7.502	ng	
90) Benzo(a)pyrene	15.12	252	507449	14.542	ng	# 97
91) Dibenzo(a,h)anthracene	16.50	278	74706	2.261	ng	# 76
92) Benzo(a,h,i)perylene	16.89	276	263073	7.916	ng	99

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

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