

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
 Data File : BF116091.D
 Acq On : 9 Aug 2019 2:04
 Operator : HP/JU
 Sample : K4154-08 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PS-8

Manual Integrations
 APPROVED

mohammad
 8/14/2019 9:21:39 AM

Quant Time: Aug 09 16:55:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	141529	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	590986	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	313781	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	509740	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	335244	20.00	ng	-0.01
87) Perylene-d12	15.46	264	364989	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	416437	49.26	ng	-0.01
7) Phenol-d6	6.47	99	575402	53.94	ng	-0.02
23) Nitrobenzene-d5	7.40	82	351432	34.33	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	136536	44.88	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	552229	32.96	ng	-0.02
79) Terphenyl-d14	12.93	244	477202	29.99	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.59	163	155790	7.292	ng	98
71) Phenanthrene	11.38	178	88456	3.394	ng	97
75) Fluoranthene	12.57	202	136528	5.004	ng	98
78) Pyrene	12.80	202	106694	4.222	ng	98
81) Benzo(a)anthracene	13.99	228	49349	2.303	ng	98
83) Chrysene	14.02	228	52248	2.512	ng	98
88) Benzo(b)fluoranthene	15.04	252	63576m	3.012	ng	

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

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