

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
 Data File : BF116740.D
 Acq On : 1 Sep 2019 14:10
 Operator : HP/JU
 Sample : K4525-19
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T9-12-13-C-(0-1.9)

Manual Integrations
 APPROVED

mohammad
 9/5/2019 9:54:14 AM

Quant Time: Sep 02 05:35:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	124804	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	507315	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	271536	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	456969	20.00	ng	-0.01
76) Chrysene-d12	13.99	240	403153	20.00	ng	-0.01
87) Perylene-d12	15.44	264	432758	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	434691	56.43	ng	-0.01
7) Phenol-d6	6.47	99	671017	66.33	ng	-0.01
23) Nitrobenzene-d5	7.40	82	407496	45.85	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	166709	91.84	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	813121	50.51	ng	-0.01
79) Terphenyl-d14	12.94	244	957713	43.95	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.59	163	42814	2.284	ng	98
71) Phenanthrene	11.38	178	58176	2.287	ng	95
75) Fluoranthene	12.56	202	113206	4.528	ng	97
78) Pyrene	12.79	202	98362	2.745	ng	99
88) Benzo(b)fluoranthene	15.02	252	62232m	2.379	ng	

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

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