

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091818\
 Data File : BF109116.D
 Acq On : 19 Sep 2018 00:32
 Operator : JU/SJ
 Sample : J4967-03 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-UST-091718-03

Manual Integrations
 APPROVED

Sohil
 9/19/2018 4:24:09 PM

Quant Time: Sep 19 05:41:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 14 13:26:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	114107	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	386010	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	159530	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	301408	20.00	ng	0.00
75) Chrysene-d12	13.85	240	345182	20.00	ng	0.00
86) Perylene-d12	15.26	264	280067	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	116977	17.03	ng	0.00
7) Phenol-d6	6.35	99	148011	16.71	ng	-0.01
23) Nitrobenzene-d5	7.28	82	87669	12.74	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	33435	19.31	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	155158	15.24	ng	0.00
78) Terphenyl-d14	12.81	244	168599	11.58	ng	0.00
Target Compounds						
70) Phenanthrene	11.25	178	37711	2.552	ng	99
74) Fluoranthene	12.44	202	80279	5.147	ng	94
77) Pyrene	12.66	202	76351	2.978	ng	98
82) Chrysene	13.88	228	45268	2.165	ng	96
87) Benzo(b)fluoranthene	14.86	252	49928m	3.222	ng	
89) Benzo(a)pyrene	15.20	252	31412	2.283	ng	# 95

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

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