

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091818\
 Data File : BF109115.D
 Acq On : 19 Sep 2018 00:05
 Operator : JU/SJ
 Sample : J4967-02 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-UST-091718-02

Manual Integrations
 APPROVED

Sohil
 9/19/2018 4:24:06 PM

Quant Time: Sep 19 05:39:10 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	116021	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	403706	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	164438	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	299136	20.00	ng	0.00
75) Chrysene-d12	13.86	240	327148	20.00	ng	0.00
86) Perylene-d12	15.27	264	255935	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	98890	14.16	ng	0.00
7) Phenol-d6	6.35	99	127652	14.17	ng	-0.01
23) Nitrobenzene-d5	7.28	82	79967	11.11	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	25611	14.35	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	145959	13.91	ng	0.00
78) Terphenyl-d14	12.81	244	147028	10.65	ng	0.00
Target Compounds						
70) Phenanthrene	11.25	178	46487	3.170	ng	97
74) Fluoranthene	12.44	202	69261	4.474	ng	95
77) Pyrene	12.66	202	60150	2.475	ng	98
82) Chrysene	13.88	228	41720	2.105	ng	# 95
87) Benzo(b)fluoranthene	14.87	252	43347m	3.061	ng	

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

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