

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092217\
 Data File : BF098715.D
 Acq On : 23 Sep 2017 00:06
 Operator : SJ/JU
 Sample : I5304-04 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2P-ENV-118-6(0-5)

Manual Integrations
 APPROVED

Sohil
 9/25/2017 6:21:19 PM

Quant Time: Sep 23 02:47:55 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 10:38:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	167911	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	683345	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	284316	20.00	ng	0.00
63) Phenanthrene-d10	11.45	188	404494	20.00	ng	0.00
75) Chrysene-d12	14.09	240	358301	20.00	ng	0.00
86) Perylene-d12	15.59	264	317556	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	233768	21.14	ng	0.00
7) Phenol-d6	6.53	99	283288	21.16	ng	-0.01
23) Nitrobenzene-d5	7.48	82	163145	16.19	ng	-0.01
41) 2,4,6-Tribromophenol	10.75	330	46286	17.80	ng	0.00
44) 2-Fluorobiphenyl	9.28	172	301705	16.89	ng	0.00
78) Terphenyl-d14	13.03	244	202559	12.14	ng	0.00
Target Compounds						
70) Phenanthrene	11.47	178	48957	2.231	ng	97
74) Fluoranthene	12.66	202	158414	7.244	ng	97
77) Pyrene	12.89	202	142138	5.343	ng	98
80) Benzo(a)anthracene	14.08	228	63228	2.930	ng	99
82) Chrysene	14.12	228	57583	2.763	ng	98
87) Benzo(b)fluoranthene	15.14	252	43008m	2.335	ng	

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

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