

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
 Data File : BF109342.D
 Acq On : 26 Sep 2018 17:45
 Operator : JU/SJ
 Sample : J5070-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-RT-4005

Manual Integrations
 APPROVED

Sohil
 9/27/2018 1:31:20 PM

Quant Time: Sep 27 09:12:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	100525	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	387824	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	184534	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	343084	20.00	ng	0.00
75) Chrysene-d12	13.84	240	244601	20.00	ng	0.00
86) Perylene-d12	15.24	264	276215	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	520060	85.21	ng	0.00
7) Phenol-d6	6.35	99	671510	86.22	ng	-0.01
23) Nitrobenzene-d5	7.26	82	418727	63.74	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	174605	95.98	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	742784	60.71	ng	-0.01
78) Terphenyl-d14	12.79	244	612476	61.45	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	270684	20.168	ng	99
74) Fluoranthene	12.42	202	54010	2.950	ng	96
77) Pyrene	12.65	202	48058	2.741	ng	99
82) Chrysene	13.86	228	32726	2.241	ng	95
87) Benzo(b)fluoranthene	14.84	252	61028m	4.021	ng	
89) Benzo(a)pyrene	15.18	252	31079	2.272	ng	97

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

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