

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109383.D
 Acq On : 27 Sep 2018 13:35
 Operator : JU/SJ
 Sample : J5066-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-Q1-HI1

Manual Integrations
 APPROVED

Sohil
 9/28/2018 2:39:25 PM

Quant Time: Sep 27 14:51:49 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	115580	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	457938	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	229288	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	433185	20.00	ng	0.00
75) Chrysene-d12	13.84	240	267743	20.00	ng	0.00
86) Perylene-d12	15.24	264	282937	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	522344	74.44	ng	0.00
7) Phenol-d6	6.35	99	681347	76.09	ng	-0.01
23) Nitrobenzene-d5	7.26	82	431456	55.62	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	162414	71.85	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	740359	48.70	ng	-0.01
78) Terphenyl-d14	12.80	244	686851	62.96	ng	0.00
Target Compounds						
10) Phenol	6.36	94	26045	2.568	ng	97
49) Dimethylphthalate	9.46	163	259434	15.556	ng	99
70) Phenanthrene	11.23	178	118578	5.482	ng	98
74) Fluoranthene	12.42	202	121018	5.236	ng	96
77) Pyrene	12.65	202	102874	5.361	ng	97
80) Benzo(a)anthracene	13.83	228	44417	2.754	ng	97
82) Chrysene	13.86	228	36857	2.306	ng	96
87) Benzo(b)fluoranthene	14.84	252	51456m	3.310	ng	
89) Benzo(a)pyrene	15.18	252	34738	2.480	ng	96

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

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