

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092317\
 Data File : BF098747.D
 Acq On : 23 Sep 2017 19:04
 Operator : SJ/JU
 Sample : I5340-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LT-TS-009

Manual Integrations
 APPROVED

Sohil
 9/25/2017 6:26:21 PM

Quant Time: Sep 25 07:13:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	150224	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	611243	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	252191	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	367369	20.00	ng	0.00
75) Chrysene-d12	14.08	240	237060	20.00	ng	0.00
86) Perylene-d12	15.57	264	245474	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1098023	116.36	ng	0.00
7) Phenol-d6	6.55	99	1334156	114.61	ng	0.00
23) Nitrobenzene-d5	7.48	82	824505	86.51	ng	0.00
41) 2,4,6-Tribromophenol	10.74	330	211017	102.26	ng	0.00
44) 2-Fluorobiphenyl	9.28	172	1302240	86.20	ng	0.00
78) Terphenyl-d14	13.03	244	806596	77.38	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.67	163	222665	11.698	ng	99
70) Phenanthrene	11.47	178	41295	2.100	ng	96
74) Fluoranthene	12.65	202	75284	3.781	ng	98
77) Pyrene	12.89	202	72903	4.033	ng	98
80) Benzo(a)anthracene	14.07	228	29158	2.031	ng	93
82) Chrysene	14.10	228	32363	2.368	ng	98
87) Benzo(b)fluoranthene	15.13	252	40422m	2.678	ng	
89) Benzo(a)pyrene	15.50	252	30172	2.178	ng	98

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

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