

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
 Data File : BF099206.D
 Acq On : 7 Oct 2017 20:49
 Operator : SJ/JU
 Sample : I5586-09 10X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G54A-0-3.5

Manual Integrations
 APPROVED

Sohil
 10/9/2017 6:34:58 PM

Quant Time: Oct 08 05:04:22 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Oct 08 03:54:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	119563	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	441915	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	170878	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	260189	20.00	ng	0.00
75) Chrysene-d12	14.03	240	212088	20.00	ng	0.00
86) Perylene-d12	15.51	264	162134	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	61085	8.13	ng	0.00
7) Phenol-d6	6.49	99	77526	8.37	ng	0.00
23) Nitrobenzene-d5	7.42	82	39686	5.76	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	6363	4.55	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	84355	8.24	ng	0.00
78) Terphenyl-d14	12.97	244	60830	6.52	ng	0.00
Target Compounds						
70) Phenanthrene	11.42	178	221652	15.915	ng	98
74) Fluoranthene	12.60	202	315323	22.359	ng	97
77) Pyrene	12.83	202	222100	13.733	ng	98
80) Benzo(a)anthracene	14.02	228	38243	2.978	ng	96
82) Chrysene	14.06	228	60687	4.964	ng	97
87) Benzo(b)fluoranthene	15.07	252	39772m	3.990	ng	

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

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