

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110917\
 Data File : BF100357.D
 Acq On : 9 Nov 2017 22:50
 Operator : SJ/JU
 Sample : I6379-07 50X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-(1-4)

Manual Integrations
 APPROVED

Sohil
 11/10/2017 1:15:22 PM

Quant Time: Nov 10 02:58:02 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 16:04:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	163585	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	657109	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	304733	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	548160	20.00	ng	0.00
75) Chrysene-d12	14.06	240	332508	20.00	ng	0.00
86) Perylene-d12	15.54	264	355973	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	18531	1.82	ng	0.00
7) Phenol-d6	6.50	99	23658	1.88	ng	-0.02
23) Nitrobenzene-d5	7.45	82	12692	1.28	ng	-0.02
41) 2,4,6-Tribromophenol	10.72	330	5381	1.73	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	29385	1.47	ng	-0.01
78) Terphenyl-d14	13.00	244	22692	1.57	ng	0.00
Target Compounds						
70) Phenanthrene	11.45	178	172686	5.983	ng	98
74) Fluoranthene	12.63	202	219654	7.181	ng	96
77) Pyrene	12.86	202	183029	7.722	ng	97
80) Benzo(a)anthracene	14.05	228	73254	3.658	ng	95
82) Chrysene	14.09	228	68585	3.537	ng	95
85) Indeno(1,2,3-cd)pyrene	17.02	276	34949	2.228	ng	# 92
87) Benzo(b)fluoranthene	15.10	252	78874m	3.740	ng	
89) Benzo(a)pyrene	15.47	252	57722	3.087	ng	97

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

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