

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110917\
 Data File : BF100354.D
 Acq On : 9 Nov 2017 21:29
 Operator : SJ/JU
 Sample : I6381-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RT-1500

Manual Integrations
 APPROVED

Sohil
 11/10/2017 1:15:16 PM

Quant Time: Nov 10 01:57:16 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	189602	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	761213	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	355847	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	664004	20.00	ng	0.00
75) Chrysene-d12	14.06	240	422952	20.00	ng	0.00
86) Perylene-d12	15.54	264	389744	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1130548	95.58	ng	0.02
7) Phenol-d6	6.53	99	1430401	98.07	ng	0.00
23) Nitrobenzene-d5	7.46	82	886171	76.97	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	371456	102.44	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1326812	56.78	ng	0.00
78) Terphenyl-d14	13.01	244	1309325	71.13	ng	0.00
Target Compounds						
10) Phenol	6.54	94	46880	3.062	ng	93
70) Phenanthrene	11.44	178	71644	2.049	ng	98
74) Fluoranthene	12.63	202	171674	4.633	ng	97
77) Pyrene	12.86	202	160180	5.313	ng	99
80) Benzo(a)anthracene	14.05	228	55992	2.198	ng	97
82) Chrysene	14.09	228	60204	2.441	ng	98
87) Benzo(b)fluoranthene	15.10	252	66746m	2.891	ng	

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

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