

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
 Data File : BF100133.D
 Acq On : 3 Nov 2017 14:40
 Operator : SJ/JU
 Sample : I6258-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BOXWOOD-DR-2-4-12-14

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:18:44 PM

Quant Time: Nov 03 15:54:44 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 15:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	86348	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	362355	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	176122	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	320966	20.00	ng	0.00
75) Chrysene-d12	13.95	240	218685	20.00	ng	0.00
86) Perylene-d12	15.41	264	152736	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	648214	117.86	ng	0.00
7) Phenol-d6	6.42	99	782862	120.04	ng	0.00
23) Nitrobenzene-d5	7.34	82	444592	78.99	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	184227	114.78	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	868362	76.07	ng	0.00
78) Terphenyl-d14	12.88	244	759460	75.89	ng	0.00
Target Compounds						
10) Phenol	6.43	94	40240	5.620	ng	93
49) Dimethylphthalate	9.52	163	80082	5.742	ng	99
74) Fluoranthene	12.52	202	72763	3.865	ng	98
77) Pyrene	12.74	202	58525	3.520	ng	96
82) Chrysene	13.98	228	28244	2.167	ng	97
87) Benzo(b)fluoranthene	15.00	252	29099m	3.017	ng	

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

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