

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110717\
 Data File : BF100280.D
 Acq On : 7 Nov 2017 13:22
 Operator : SJ/JU
 Sample : I6336-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CF-1B(4-5)

Manual Integrations
 APPROVED

Sohil
 11/8/2017 3:54:05 PM

Quant Time: Nov 07 23:54:49 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 02:59:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	95957	20.00	ng	-0.05
21) Naphthalene-d8	7.99	136	400510	20.00	ng	-0.05
38) Acenaphthene-d10	9.74	164	191989	20.00	ng	-0.06
63) Phenanthrene-d10	11.23	188	358884	20.00	ng	-0.06
75) Chrysene-d12	13.89	240	240471	20.00	ng	-0.06
86) Perylene-d12	15.33	264	158363	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	106776	17.47	ng	-0.01
7) Phenol-d6	6.39	99	74070	10.22	ng	-0.03
23) Nitrobenzene-d5	7.29	82	238084	38.27	ng	-0.05
41) 2,4,6-Tribromophenol	10.54	330	72988	41.72	ng	-0.06
44) 2-Fluorobiphenyl	9.06	172	525328	42.22	ng	-0.05
78) Terphenyl-d14	12.82	244	448574	40.77	ng	-0.05
Target Compounds						
10) Phenol	6.41	94	30174	3.792	ng	97
49) Dimethylphthalate	9.47	163	96211	6.328	ng	100
51) Acenaphthene	9.77	154	61100	5.147	ng	98
54) Dibenzofuran	9.96	168	64570	3.854	ng	97
57) Fluorene	10.30	166	51597	3.719	ng	99
70) Phenanthrene	11.26	178	54999	2.716	ng	98

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

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