

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
 Data File : BF110410.D
 Acq On : 2 Nov 2018 14:23
 Operator : JU/SJ
 Sample : J5794-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-11-S

Manual Integrations
 APPROVED

Sohil
 11/5/2018 10:32:39 AM

Quant Time: Nov 02 23:01:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	90023	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	354060	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	147636	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	231386	20.00	ng	0.00
76) Chrysene-d12	14.13	240	210346	20.00	ng	0.00
87) Perylene-d12	15.66	264	182876	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	572426	103.55	ng	0.00
7) Phenol-d6	6.58	99	720804	101.94	ng	0.00
23) Nitrobenzene-d5	7.52	82	454725	76.18	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	133382	91.21	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	781180	83.09	ng	0.00
79) Terphenyl-d14	13.07	244	610485	60.36	ng	0.00
Target Compounds						
10) Phenol	6.59	94	41322	5.467	ng	Qvalue # 66
50) Dimethylphthalate	9.70	163	122332	11.225	ng	98
75) Fluoranthene	12.70	202	37068	3.017	ng	99
78) Pyrene	12.93	202	32920	2.183	ng	98
88) Benzo(b)fluoranthene	15.20	252	24171m	2.289	ng	

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

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