

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110718\
 Data File : BF110558.D
 Acq On : 7 Nov 2018 16:33
 Operator : JU/SJ
 Sample : J5839-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-4

Quant Time: Nov 09 10:29:06 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	6.95	152	72447	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	292992	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	140644	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	249924	20.00	ng	0.00
76) Chrysene-d12	14.13	240	159455	20.00	ng	0.00
87) Perylene-d12	15.66	264	162434	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	551256	123.91	ng	0.00
7) Phenol-d6	6.58	99	694973	122.13	ng	0.00
23) Nitrobenzene-d5	7.52	82	447749	90.64	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	168734	121.11	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	870197	97.16	ng	0.00
79) Terphenyl-d14	13.07	244	723577	94.37	ng	0.00
Target Compounds						
10) Phenol	6.59	94	46692	7.676	ng	Qvalue # 63
50) Dimethylphthalate	9.70	163	138111	13.303	ng	99
71) Phenanthrene	11.51	178	67968	5.197	ng	99
75) Fluoranthene	12.70	202	56055	4.224	ng	98
78) Pyrene	12.94	202	43669	3.820	ng	97
81) Benzo(a)anthracene	14.13	228	18919	2.001	ng	96

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

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