

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061318\
 Data File : BF106392.D
 Acq On : 13 Jun 2018 15:44
 Operator : JU/SJ
 Sample : J3459-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-RO-341

Manual Integrations
 APPROVED

Sohil
 6/14/2018 2:47:38 PM

Quant Time: Jun 14 03:07:38 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 10:51:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	135078	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	515706	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	254243	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	510807	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	356259	20.00	ng	-0.02
86) Perylene-d12	15.68	264	378451	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	591378	74.10	ng	-0.02
7) Phenol-d6	6.61	99	760321	75.40	ng	-0.02
23) Nitrobenzene-d5	7.52	82	461817	52.68	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	209072	85.47	ng	-0.01
44) 2-Fluorobiphenyl	9.32	172	841469	50.88	ng	-0.01
78) Terphenyl-d14	13.08	244	906919	63.23	ng	0.00
Target Compounds				Qvalue		
31) Naphthalene	8.27	128	138224	5.929	ng	99
37) 2-Methylnaphthalene	8.96	142	91520	5.764	ng	98
45) 1,1'-Biphenyl	9.42	154	58022	2.946	ng	95
49) Dimethylphthalate	9.71	163	84467	4.650	ng	99
51) Acenaphthene	10.04	154	257791	16.549	ng	97
54) Dibenzofuran	10.21	168	299870	14.295	ng	99
57) Fluorene	10.55	166	244538	14.714	ng	99
70) Phenanthrene	11.52	178	627608	25.088	ng	98
71) Anthracene	11.57	178	56397	2.250	ng	98
72) Carbazole	11.72	167	210289	9.089	ng	98
74) Fluoranthene	12.71	202	341696	12.808	ng	96
77) Pvrene	12.94	202	237430	10.645	ng	99
80) Benzo(a)anthracene	14.14	228	48622	2.293	ng	93
82) Chrysene	14.17	228	51696	2.671	ng	95
87) Benzo(b)fluoranthene	15.22	252	53720m	2.348	ng	

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

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