

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
 Data File : BF109346.D
 Acq On : 26 Sep 2018 19:29
 Operator : JU/SJ
 Sample : J4773-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-092518-A

Manual Integrations
 APPROVED

Sohil
 9/27/2018 1:31:31 PM

Quant Time: Sep 27 06:27:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	107151	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	417427	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	192749	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	311205	20.00	ng	0.00
75) Chrysene-d12	13.84	240	279902	20.00	ng	0.00
86) Perylene-d12	15.24	264	244088	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	405174	62.28	ng	0.00
7) Phenol-d6	6.34	99	533779	64.30	ng	-0.02
23) Nitrobenzene-d5	7.26	82	327705	46.34	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	119445	62.86	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	590997	46.24	ng	-0.01
78) Terphenyl-d14	12.79	244	446447	39.14	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	201240	14.355	ng	98
70) Phenanthrene	11.23	178	59073	3.802	ng	98
74) Fluoranthene	12.42	202	119575	7.201	ng	95
77) Pyrene	12.65	202	126425	6.302	ng	99
80) Benzo(a)anthracene	13.83	228	68000	4.033	ng	97
82) Chrysene	13.86	228	61826	3.700	ng	95
87) Benzo(b)fluoranthene	14.84	252	70367m	5.246	ng	
89) Benzo(a)pyrene	15.18	252	48875	4.044	ng	96
91) Benzo(g,h,i)perylene	16.99	276	22156	2.274	ng	91

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

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