

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
 Data File : BF109349.D
 Acq On : 26 Sep 2018 20:48
 Operator : JU/SJ
 Sample : J4773-17
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 9/27/2018 1:31:36 PM

Quant Time: Sep 27 09:39:36 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	6.71	152	88215	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	288495	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	120644	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	248023	20.00	ng	0.00
75) Chrysene-d12	13.84	240	268458	20.00	ng	0.00
86) Perylene-d12	15.24	264	210238	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	416853	77.83	ng	0.00
7) Phenol-d6	6.35	99	505957	74.03	ng	-0.01
23) Nitrobenzene-d5	7.26	82	304028	62.21	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	114100	95.94	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	475671	59.46	ng	-0.01
78) Terphenyl-d14	12.80	244	530860	48.53	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	154821	17.644	ng	98
70) Phenanthrene	11.24	178	38739	3.128	ng	97
74) Fluoranthene	12.42	202	57635	4.355	ng	96
77) Pyrene	12.64	202	71468	3.715	ng	98
80) Benzo(a)anthracene	13.83	228	34155	2.112	ng	95
82) Chrysene	13.87	228	33916	2.116	ng	95
87) Benzo(b)fluoranthene	14.84	252	28736m	2.487	ng	
89) Benzo(a)pyrene	15.18	252	21127	2.030	ng	# 91

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

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