

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091018\
 Data File : BF108892.D
 Acq On : 10 Sep 2018 14:56
 Operator : JU/SJ
 Sample : J4776-03 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-090718-B

Manual Integrations
 APPROVED

Sohil
 9/11/2018 12:01:32 PM

Quant Time: Sep 11 09:14:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	212640	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	794104	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	344450	20.00	ng	0.00
63) Phenanthrene-d10	11.25	188	540285	20.00	ng	0.00
75) Chrysene-d12	13.87	240	509122	20.00	ng	0.00
86) Perylene-d12	15.28	264	445817	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	256246	19.96	ng	0.00
7) Phenol-d6	6.36	99	329550	19.95	ng	0.00
23) Nitrobenzene-d5	7.29	82	190415	15.01	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	57142	17.38	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	368298	18.24	ng	0.00
78) Terphenyl-d14	12.83	244	272243	12.47	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	84770	3.519	ng	Qvalue # 98
70) Phenanthrene	11.27	178	126111	4.888	ng	99
71) Anthracene	11.32	178	50512	2.045	ng	98
74) Fluoranthene	12.45	202	188464	7.378	ng	97
77) Pyrene	12.68	202	249317	6.364	ng	100
80) Benzo(a)anthracene	13.87	228	125871	3.857	ng	93
82) Chrysene	13.90	228	129477	4.099	ng	97
87) Benzo(b)fluoranthene	14.88	252	147480m	6.084	ng	
88) Benzo(k)fluoranthene	14.91	252	44649m	2.000	ng	
89) Benzo(a)pyrene	15.22	252	108796	4.974	ng	# 93
91) Benzo(g,h,i)perylene	17.05	276	50221	2.602	ng	# 92

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

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