

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
 Data File : BF109353.D
 Acq On : 26 Sep 2018 22:33
 Operator : JU/SJ
 Sample : J5081-17
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-S

Manual Integrations
 APPROVED

Sohil
 9/28/2018 5:01:58 PM

Quant Time: Sep 27 09:44:35 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	90726	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	317453	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	137410	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	260546	20.00	ng	0.00
75) Chrysene-d12	13.84	240	277089	20.00	ng	0.00
86) Perylene-d12	15.24	264	229813	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	474893	86.21	ng	0.00
7) Phenol-d6	6.35	99	583149	82.97	ng	-0.01
23) Nitrobenzene-d5	7.26	82	364623	67.80	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	132785	98.03	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	605646	66.47	ng	-0.01
78) Terphenyl-d14	12.80	244	628414	55.66	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	206115	20.623	ng	99
74) Fluoranthene	12.42	202	88032	6.332	ng	94
77) Pyrene	12.64	202	126105	6.350	ng	97
80) Benzo(a)anthracene	13.83	228	92683m	5.553	ng	
82) Chrysene	13.86	228	90724	5.484	ng	98
85) Indeno(1,2,3-cd)pyrene	16.59	276	35387	2.639	ng	93
87) Benzo(b)fluoranthene	14.84	252	91241m	7.225	ng	
88) Benzo(k)fluoranthene	14.87	252	27484m	2.294	ng	
89) Benzo(a)pyrene	15.18	252	67448	5.927	ng	98
91) Benzo(g,h,i)perylene	16.99	276	37263	4.061	ng	93

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

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