

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109382.D
 Acq On : 27 Sep 2018 13:09
 Operator : JU/SJ
 Sample : J5081-19
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-S

Manual Integrations
 APPROVED

Sohil
 9/28/2018 2:38:46 PM

Quant Time: Sep 27 14:21:48 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	112364	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	446340	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	225413	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	442150	20.00	ng	0.00
75) Chrysene-d12	13.84	240	285127	20.00	ng	0.00
86) Perylene-d12	15.24	264	279942	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	478548	70.15	ng	0.00
7) Phenol-d6	6.35	99	620447	71.27	ng	-0.01
23) Nitrobenzene-d5	7.26	82	403590	53.38	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	161113	72.50	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	774034	51.79	ng	-0.01
78) Terphenyl-d14	12.80	244	722907	62.22	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	248648	15.166	ng	99
74) Fluoranthene	12.42	202	73684	3.123	ng	95
77) Pyrene	12.64	202	74091	3.626	ng	96
80) Benzo(a)anthracene	13.83	228	46920m	2.732	ng	
82) Chrysene	13.86	228	42335	2.487	ng	97
87) Benzo(b)fluoranthene	14.84	252	55389m	3.601	ng	
89) Benzo(a)pyrene	15.18	252	43084	3.108	ng	98
91) Benzo(g,h,i)perylene	16.99	276	22815	2.041	ng	94

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

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