

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092818\
 Data File : BF109437.D
 Acq On : 28 Sep 2018 18:32
 Operator : JU/SJ
 Sample : J5077-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PARCEL-190-092418-1

Manual Integrations
 APPROVED

Sohil
 10/1/2018 9:57:09 AM

Quant Time: Sep 29 02:22:59 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	99506	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	329610	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	129790	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	241265	20.00	ng	0.00
75) Chrysene-d12	13.84	240	191376	20.00	ng	0.00
86) Perylene-d12	15.25	264	141289	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	439941	72.82	ng	0.00
7) Phenol-d6	6.35	99	535479	69.46	ng	-0.01
23) Nitrobenzene-d5	7.27	82	322883	57.83	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	99023	77.39	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	461634	53.64	ng	-0.01
78) Terphenyl-d14	12.80	244	412767	52.93	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	181746	19.253	ng	99
74) Fluoranthene	12.42	202	66288	5.149	ng	98
77) Pyrene	12.65	202	55367	4.037	ng	98
80) Benzo(a)anthracene	13.83	228	26943	2.337	ng	98
82) Chrysene	13.87	228	33528	2.935	ng	94
87) Benzo(b)fluoranthene	14.85	252	39757m	5.121	ng	
89) Benzo(a)pyrene	15.19	252	22165	3.168	ng	# 80
91) Benzo(g,h,i)perylene	17.01	276	17365	3.079	ng	# 88

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

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