

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109401.D
 Acq On : 27 Sep 2018 21:28
 Operator : JU/SJ
 Sample : J5077-06
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PARCEL-193-092418-2

Manual Integrations
 APPROVED

Sohil
 9/28/2018 2:40:24 PM

Quant Time: Sep 28 05:43:57 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	109562	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	407663	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	171378	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	270430	20.00	ng	0.00
75) Chrysene-d12	13.84	240	254925	20.00	ng	0.00
86) Perylene-d12	15.24	264	190862	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	534457	80.35	ng	0.00
7) Phenol-d6	6.35	99	690907	81.40	ng	-0.01
23) Nitrobenzene-d5	7.27	82	432381	62.61	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	131945	78.10	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	736676	64.83	ng	-0.01
78) Terphenyl-d14	12.79	244	584182	56.24	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	267138	21.431	ng	99
70) Phenanthrene	11.23	178	57158	4.233	ng	99
74) Fluoranthene	12.42	202	106067	7.351	ng	96
77) Pyrene	12.65	202	116610	6.383	ng	99
80) Benzo(a)anthracene	13.83	228	56639	3.689	ng	92
82) Chrysene	13.86	228	60133	3.951	ng	97
87) Benzo(b)fluoranthene	14.84	252	63697m	6.073	ng	
89) Benzo(a)pyrene	15.18	252	37683	3.987	ng	# 94
91) Benzo(g,h,i)perylene	16.99	276	24541	3.221	ng	# 91

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

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