

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092218\
 Data File : BF109249.D
 Acq On : 23 Sep 2018 4:39
 Operator : JU/SJ
 Sample : J5017-04
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-091918-S-SW-058

Manual Integrations
 APPROVED

Sohil
 9/24/2018 11:39:20 AM

Quant Time: Sep 24 08:29:14 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	107727	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	381985	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	149254	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	236165	20.00	ng	0.00
75) Chrysene-d12	13.84	240	207869	20.00	ng	0.00
86) Perylene-d12	15.24	264	172745	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	729686	111.56	ng	0.00
7) Phenol-d6	6.35	99	952261	114.10	ng	-0.01
23) Nitrobenzene-d5	7.27	82	595311	92.00	ng	-0.01
41) 2,4,6-Tribromophenol	10.53	330	157998	107.38	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	952450	96.24	ng	0.00
78) Terphenyl-d14	12.80	244	684937	80.87	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	212236	19.551	ng	# 98
70) Phenanthrene	11.24	178	54695	4.638	ng	# 90
74) Fluoranthene	12.42	202	47637	3.780	ng	93
77) Pyrene	12.65	202	64507	4.330	ng	95
80) Benzo(a)anthracene	13.83	228	28674	2.290	ng	# 83
82) Chrysene	13.86	228	26416	2.129	ng	# 83
87) Benzo(b)fluoranthene	14.84	252	22645m	2.386	ng	

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

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