

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091718\
 Data File : BF109040.D
 Acq On : 17 Sep 2018 11:36
 Operator : JU/SJ
 Sample : J4941-04 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RB47198RT

Manual Integrations
 APPROVED

Sohil
 9/18/2018 5:59:53 PM

Quant Time: Sep 17 12:02:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 14 13:26:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	86979	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	358720	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	163025	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	248411	20.00	ng	0.00
75) Chrysene-d12	13.85	240	197377	20.00	ng	0.00
86) Perylene-d12	15.25	264	229883	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	178630	34.11	ng	0.00
7) Phenol-d6	6.34	99	232292	34.40	ng	-0.02
23) Nitrobenzene-d5	7.27	82	161439	25.24	ng	-0.02
41) 2,4,6-Tribromophenol	10.53	330	45794	25.88	ng	-0.01
44) 2-Fluorobiphenyl	9.07	172	285274	27.42	ng	-0.01
78) Terphenyl-d14	12.81	244	181861	21.84	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.47	163	41702	3.579	ng	98
74) Fluoranthene	12.44	202	37498	2.917	ng	91
77) Pyrene	12.66	202	32982	2.250	ng	# 94
83) Bis(2-ethylhexyl)phthalate	13.85	149	74579	9.406	ng	97
87) Benzo(b)fluoranthene	14.85	252	28422m	2.235	ng	
89) Benzo(a)pyrene	15.19	252	22638	2.004	ng	# 88

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

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