

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
 Data File : BF109276.D
 Acq On : 24 Sep 2018 19:03
 Operator : JU/SJ
 Sample : J4770-15 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-092118-H

Manual Integrations
 APPROVED

Sohil
 9/25/2018 9:57:24 AM

Quant Time: Sep 25 05:25:02 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	96182	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	342438	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	153111	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	316331	20.00	ng	0.00
75) Chrysene-d12	13.84	240	256871	20.00	ng	0.00
86) Perylene-d12	15.25	264	174897	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	194033	33.23	ng	0.00
7) Phenol-d6	6.34	99	235680	31.63	ng	-0.02
23) Nitrobenzene-d5	7.26	82	132190	22.79	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	56204	37.24	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	237966	23.44	ng	-0.01
78) Terphenyl-d14	12.80	244	273274	26.11	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	66785	5.997	ng	Qvalue # 98
70) Phenanthrene	11.24	178	74894	4.742	ng	98
74) Fluoranthene	12.43	202	248963	14.750	ng	93
77) Pyrene	12.65	202	260171	14.132	ng	99
80) Benzo(a)anthracene	13.83	228	114979	7.431	ng	99
82) Chrysene	13.87	228	94611	6.170	ng	96
85) Indeno(1,2,3-cd)pyrene	16.60	276	44529	3.582	ng	# 89
87) Benzo(b)fluoranthene	14.85	252	97377m	10.132	ng	
88) Benzo(k)fluoranthene	14.87	252	32268m	3.538	ng	
89) Benzo(a)pyrene	15.19	252	68660	7.929	ng	96
91) Benzo(g,h,i)perylene	17.00	276	42736	6.121	ng	92

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

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