

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF063018\
 Data File : BF107025.D
 Acq On : 30 Jun 2018 19:57
 Operator : JU/SJ
 Sample : J3629-15 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A0C22-01

Manual Integrations
 APPROVED

Sohil
 7/2/2018 12:30:16 PM

Quant Time: Jul 02 07:11:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	71334	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	251512	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	109069	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	222363	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	193784	20.00	ng	-0.02
86) Perylene-d12	15.68	264	134671	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	219431	52.09	ng	0.00
7) Phenol-d6	6.60	99	262675	49.93	ng	-0.02
23) Nitrobenzene-d5	7.51	82	141071	31.17	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	64904	51.34	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	233327	25.96	ng	-0.02
78) Terphenyl-d14	13.07	244	283052	35.38	ng	-0.02
Target Compounds						
9) Benzaldehyde	6.73	77	4658	1.096	ng	86
49) Dimethylphthalate	9.70	163	27560	3.369	ng	97
66) 4-Bromophenyl-phenylether	10.79	248	4050	1.526	ng	# 1
70) Phenanthrene	11.51	178	26742	2.493	ng	99
74) Fluoranthene	12.71	202	56588	4.787	ng	94
77) Pyrene	12.94	202	64276	5.030	ng	99
80) Benzo(a)anthracene	14.13	228	25484	2.158	ng	# 90
82) Chrysene	14.17	228	26768m	2.464	ng	
83) Bis(2-ethylhexyl)phthalate	14.09	149	20034	2.983	ng	# 95
85) Indeno(1,2,3-cd)pyrene	17.25	276	11476	1.245	ng	92
87) Benzo(b)fluoranthene	15.22	252	22320m	2.668	ng	
89) Benzo(a)pyrene	15.61	252	16018m	2.154	ng	
91) Benzo(a,h,i)perylene	17.74	276	13134m	2.132	ng	

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

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