

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070918\
 Data File : BF107238.D
 Acq On : 9 Jul 2018 19:28
 Operator : JU/SJ
 Sample : J3879-09RE
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SURCHARGE-SP-5RE

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:46:45 PM

Quant Time: Jul 10 04:32:44 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	46238	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	167923	20.00	ng	-0.02
38) Acenaphthene-d10	10.00	164	77511	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	144359	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	108428	20.00	ng	-0.01
86) Perylene-d12	15.69	264	79640	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	289410	105.99	ng	0.00
7) Phenol-d6	6.60	99	357692	104.89	ng	-0.01
23) Nitrobenzene-d5	7.51	82	220254	72.89	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	100579	111.96	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	386056	84.16	ng	-0.02
78) Terphenyl-d14	13.08	244	428824	95.80	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.70	163	111089	19.109	ng	Qvalue 98
57) Fluorene	10.54	166	11441	2.178	ng	# 96
70) Phenanthrene	11.51	178	153228	22.007	ng	98
71) Anthracene	11.57	178	54712	7.698	ng	99
74) Fluoranthene	12.72	202	340262	44.338	ng	94
77) Pyrene	12.95	202	305625	42.744	ng	99
80) Benzo(a)anthracene	14.14	228	155539m	23.537	ng	
82) Chrysene	14.18	228	148110m	24.362	ng	
85) Indeno(1,2,3-cd)pyrene	17.29	276	57633	11.171	ng	# 86
87) Benzo(b)fluoranthene	15.24	252	144869m	29.283	ng	
88) Benzo(k)fluoranthene	15.27	252	50584m	10.737	ng	
89) Benzo(a)pyrene	15.63	252	98585	22.416	ng	97
90) Dibenzo(a,h)anthracene	17.29	278	15632	4.194	ng	# 88
91) Benzo(a,h,i)perylene	17.79	276	56069	15.391	ng	# 86

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

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