

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102918\
 Data File : BF110268.D
 Acq On : 29 Oct 2018 18:10
 Operator : JU/SJ
 Sample : J5696-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DL-SED-02

Manual Integrations
 APPROVED

Sohil
 10/30/2018 2:24:11 PM

Quant Time: Oct 30 07:53:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 17:29:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	127209	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	510010	20.00	ng	0.00
39) Acenaphthene-d10	10.01	164	222909	20.00	ng	0.00
64) Phenanthrene-d10	11.50	188	351497	20.00	ng	0.00
76) Chrysene-d12	14.16	240	241424	20.00	ng	0.00
87) Perylene-d12	15.69	264	200634	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	755708	96.74	ng	0.00
7) Phenol-d6	6.59	99	963222	96.40	ng	0.00
23) Nitrobenzene-d5	7.53	82	601162	69.91	ng	0.00
42) 2,4,6-Tribromophenol	10.80	330	169054	76.56	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	843790	59.44	ng	0.00
79) Terphenyl-d14	13.09	244	508316	43.79	ng	0.00
Target Compounds						
10) Phenol	6.60	94	62612	5.862	ng	Qvalue # 66
50) Dimethylphthalate	9.72	163	186699	11.346	ng	100
71) Phenanthrene	11.53	178	270935	14.731	ng	99
72) Anthracene	11.58	178	56294	3.054	ng	100
75) Fluoranthene	12.72	202	397141	21.276	ng	100
78) Pyrene	12.95	202	337972	19.527	ng	98
81) Benzo(a)anthracene	14.14	228	154684	10.804	ng	100
83) Chrysene	14.18	228	165201	12.149	ng	98
84) Bis(2-ethylhexyl)phthalate	14.11	149	22361	2.202	ng	99
86) Indeno(1,2,3-cd)pyrene	17.25	276	66547	5.899	ng	98
88) Benzo(b)fluoranthene	15.23	252	207692m	17.926	ng	
89) Benzo(k)fluoranthene	15.26	252	51656m	4.535	ng	
90) Benzo(a)pyrene	15.62	252	119040	11.166	ng	98
92) Benzo(a,h,i)perylene	17.73	276	68800	7.590	ng	97

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

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