

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122818\
 Data File : BF111792.D
 Acq On : 28 Dec 2018 16:50
 Operator : JU/SJ
 Sample : J6193-13 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-122718-B

Manual Integrations
 APPROVED

Sohil
 1/2/2019 2:51:36 PM

Quant Time: Dec 29 01:41:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	152479	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	533106	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	213611	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	341920	20.00	ng	0.01
76) Chrysene-d12	14.07	240	386847	20.00	ng	0.02
87) Perylene-d12	15.56	264	316356	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	463655	51.83	ng	0.02
7) Phenol-d6	6.54	99	568490	49.94	ng	0.02
23) Nitrobenzene-d5	7.46	82	350438	38.26	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	115590	45.80	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	577662	39.42	ng	0.00
79) Terphenyl-d14	13.01	244	519109	25.45	ng	0.01
Target Compounds						
10) Phenol	6.55	94	38023	2.960	ng	91
50) Dimethylphthalate	9.65	163	74453	4.766	ng	99
71) Phenanthrene	11.45	178	41400	2.283	ng	98
75) Fluoranthene	12.64	202	70166	3.821	ng	94
78) Pyrene	12.87	202	72904	2.669	ng	95
88) Benzo(b)fluoranthene	15.12	252	37204m	2.012	ng	

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

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