

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122118\
 Data File : BF111684.D
 Acq On : 21 Dec 2018 13:44
 Operator : JU/SJ
 Sample : J6362-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-3

Manual Integrations
 APPROVED

Sohil
 12/26/2018 3:10:40 PM

Quant Time: Dec 21 23:09:14 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	6.89	152	208171	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	772768	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	357942	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	554186	20.00	ng	0.00
76) Chrysene-d12	14.05	240	461284	20.00	ng	0.00
87) Perylene-d12	15.53	264	342172	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1245505	101.98	ng	0.00
7) Phenol-d6	6.53	99	1629974	104.87	ng	0.00
23) Nitrobenzene-d5	7.45	82	1041760	78.47	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	382841	90.52	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1862636	75.85	ng	0.00
79) Terphenyl-d14	13.00	244	1467830	60.34	ng	0.00
Target Compounds						
10) Phenol	6.53	94	114417	6.524	ng	95
50) Dimethylphthalate	9.64	163	171242	6.542	ng	97
75) Fluoranthene	12.62	202	75228	2.528	ng	96
78) Pyrene	12.85	202	84251	2.586	ng	97
88) Benzo(b)fluoranthene	15.09	252	58009m	2.901	ng	
90) Benzo(a)pyrene	15.46	252	37937	2.088	ng	98

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

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