

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111534.D
 Acq On : 17 Dec 2018 12:48
 Operator : JU/SJ
 Sample : J6386-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS005-0006-01

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:47:43 PM

Quant Time: Dec 18 04:34:41 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	152169	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	593078	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	231515	20.00	ng	-0.01
64) Phenanthrene-d10	11.31	188	339019	20.00	ng	0.00
76) Chrysene-d12	13.94	240	331255	20.00	ng	0.00
87) Perylene-d12	15.39	264	338675	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	683894	74.79	ng	0.00
7) Phenol-d6	6.43	99	859464	70.66	ng	0.00
23) Nitrobenzene-d5	7.35	82	523524	52.57	ng	-0.01
42) 2,4,6-Tribromophenol	10.61	330	113001	54.49	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	817169	54.52	ng	-0.01
79) Terphenyl-d14	12.89	244	538687	38.19	ng	0.00
Target Compounds						
10) Phenol	6.45	94	51315	3.787	ng	97
50) Dimethylphthalate	9.54	163	86539	5.190	ng	96
80) Butylbenzylphthalate	13.37	149	99570	8.747	ng	89
86) Indeno(1,2,3-cd)pyrene	16.79	276	35350	2.581	ng	96
88) Benzo(b)fluoranthene	14.97	252	69396m	3.515	ng	
90) Benzo(a)pyrene	15.32	252	40828	2.293	ng	99
92) Benzo(g,h,i)perylene	17.22	276	33843	2.455	ng	95

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

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