

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041119\
 Data File : BF113728.D
 Acq On : 11 Apr 2019 19:00
 Operator : JU/SJ
 Sample : K2343-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7-S

Manual Integrations
 APPROVED

Sohil
 4/12/2019 2:29:02 PM

Quant Time: Apr 12 02:16:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	120816	20.00	ng	-0.02
21) Naphthalene-d8	7.96	136	467710	20.00	ng	-0.01
39) Acenaphthene-d10	9.70	164	224547	20.00	ng	-0.02
64) Phenanthrene-d10	11.19	188	455593	20.00	ng	-0.02
76) Chrysene-d12	13.82	240	341667	20.00	ng	-0.02
87) Perylene-d12	15.23	264	327079	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	798796	112.74	ng	-0.01
7) Phenol-d6	6.33	99	1041895	119.31	ng	-0.01
23) Nitrobenzene-d5	7.24	82	608546	76.21	ng	-0.02
42) 2,4,6-Tribromophenol	10.50	330	265323	98.36	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	1174343	84.19	ng	-0.02
79) Terphenyl-d14	12.76	244	1217577	72.56	ng	-0.03
Target Compounds						
50) Dimethylphthalate	9.43	163	106150	6.413	ng	100
71) Phenanthrene	11.21	178	102943	4.547	ng	97
75) Fluoranthene	12.40	202	137731	5.678	ng	99
78) Pyrene	12.63	202	120789	4.646	ng	97
81) Benzo(a)anthracene	13.81	228	54974	2.789	ng	99
83) Chrysene	13.85	228	48333	2.419	ng	99
86) Indeno(1,2,3-cd)pyrene	16.60	276	40255	2.182	ng	96
88) Benzo(b)fluoranthene	14.84	252	69165m	3.455	ng	
90) Benzo(a)pyrene	15.17	252	50521	2.840	ng	99
92) Benzo(g,h,i)perylene	17.01	276	38957	2.294	ng	95

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

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