

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041819\
 Data File : BF113819.D
 Acq On : 18 Apr 2019 16:09
 Operator : JU/SJ
 Sample : K2411-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-14-S

Manual Integrations
 APPROVED

Sohil
 4/23/2019 8:16:40 AM

Quant Time: Apr 18 16:40:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	187092	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	699519	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	341368	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	595320	20.00	ng	0.00
76) Chrysene-d12	14.04	240	437552	20.00	ng	0.00
87) Perylene-d12	15.51	264	561951	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	981240	87.21	ng	0.00
7) Phenol-d6	6.52	99	1265708	90.20	ng	0.00
23) Nitrobenzene-d5	7.45	82	766685	66.74	ng	-0.01
42) 2,4,6-Tribromophenol	10.71	330	305513	81.01	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1443699	64.50	ng	0.00
79) Terphenyl-d14	12.99	244	1413229	67.73	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	186864	7.822	ng	99
75) Fluoranthene	12.62	202	87975	2.738	ng	99
78) Pyrene	12.84	202	87175	2.827	ng	97
81) Benzo(a)anthracene	14.03	228	53515	2.092	ng	95
86) Indeno(1,2,3-cd)pyrene	16.97	276	46894	2.263	ng	# 94
88) Benzo(b)fluoranthene	15.07	252	72511m	2.070	ng	

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

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