

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041519\
 Data File : BF113763.D
 Acq On : 15 Apr 2019 15:12
 Operator : JU/SJ
 Sample : K2407-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-11-S

Manual Integrations
 APPROVED

Sohil
 4/16/2019 5:02:52 PM

Quant Time: Apr 15 23:47:20 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	123215	20.00	ng	-0.02
21) Naphthalene-d8	7.95	136	457539	20.00	ng	-0.02
39) Acenaphthene-d10	9.70	164	214651	20.00	ng	-0.02
64) Phenanthrene-d10	11.19	188	431803	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	293390	20.00	ng	-0.02
87) Perylene-d12	15.23	264	308923	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	679292	94.01	ng	-0.01
7) Phenol-d6	6.33	99	912626	102.48	ng	-0.01
23) Nitrobenzene-d5	7.24	82	516759	66.16	ng	-0.02
42) 2,4,6-Tribromophenol	10.50	330	223996	86.87	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	971780	72.88	ng	-0.02
79) Terphenyl-d14	12.77	244	1056584	73.32	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.43	163	97245	6.146	ng	Qvalue # 99
71) Phenanthrene	11.21	178	148890	6.939	ng	98
75) Fluoranthene	12.40	202	263116	11.444	ng	97
78) Pyrene	12.63	202	236500	10.594	ng	99
81) Benzo(a)anthracene	13.82	228	111188	6.570	ng	96
83) Chrysene	13.85	228	93986	5.478	ng	99
86) Indeno(1,2,3-cd)pyrene	16.61	276	82651	5.218	ng	94
88) Benzo(b)fluoranthene	14.84	252	131730m	6.966	ng	
89) Benzo(k)fluoranthene	14.87	252	51179m	3.016	ng	
90) Benzo(a)pyrene	15.18	252	104785	6.236	ng	99
92) Benzo(g,h,i)perylene	17.02	276	87212	5.438	ng	96

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

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