

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
 Data File : BF112077.D
 Acq On : 17 Jan 2019 9:32
 Operator : JU/SJ
 Sample : K1147-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2

Manual Integrations
 APPROVED

Sohil
 1/17/2019 2:55:37 PM

Quant Time: Jan 17 10:29:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 16 18:22:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	280812	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1091361	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	532379	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	964695	20.00	ng	0.00
76) Chrysene-d12	14.03	240	736628	20.00	ng	0.00
87) Perylene-d12	15.50	264	695320	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1527478	98.73	ng	0.00
7) Phenol-d6	6.50	99	1986533	102.63	ng	0.00
23) Nitrobenzene-d5	7.42	82	1215686	76.96	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	604797	105.31	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2479501	68.28	ng	0.00
79) Terphenyl-d14	12.97	244	2390741	69.04	ng	0.00
Target Compounds						
10) Phenol	6.52	94	44165	2.088	ng	83
50) Dimethylphthalate	9.61	163	655573	16.885	ng	# 97
75) Fluoranthene	12.60	202	234782	4.580	ng	96
78) Pyrene	12.83	202	246228	5.066	ng	96
81) Benzo(a)anthracene	14.02	228	108612	2.572	ng	96
83) Chrysene	14.05	228	142022m	3.554	ng	
86) Indeno(1,2,3-cd)pyrene	16.97	276	91729	2.488	ng	96
88) Benzo(b)fluoranthene	15.07	252	212942m	5.395	ng	
90) Benzo(a)pyrene	15.43	252	109241	3.018	ng	# 97
92) Benzo(g,h,i)perylene	17.42	276	105179	3.321	ng	99

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

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