

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100119\
 Data File : BF117288.D
 Acq On : 1 Oct 2019 20:29
 Operator : JU
 Sample : K4659-04 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-093019-B

Manual Integrations
 APPROVED

mohammad
 10/2/2019 3:56:51 PM

Quant Time: Oct 02 06:30:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	50084	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	191258	20.00	ng	-0.02
39) Acenaphthene-d10	9.85	164	82791	20.00	ng	-0.01
64) Phenanthrene-d10	11.33	188	107722	20.00	ng	0.00
76) Chrysene-d12	13.97	240	79729	20.00	ng	0.00
87) Perylene-d12	15.43	264	66180	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	119579	37.28	ng	0.00
7) Phenol-d6	6.45	99	163580	39.46	ng	-0.01
23) Nitrobenzene-d5	7.37	82	96697	26.56	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	22320	32.26	ng	-0.01
45) 2-Fluorobiphenyl	9.16	172	169270	31.97	ng	-0.01
79) Terphenyl-d14	12.92	244	119786	28.08	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	36573	6.239	ng	# 97
71) Phenanthrene	11.35	178	14613	2.388	ng	# 96
75) Fluoranthene	12.54	202	22331	3.552	ng	96
78) Pyrene	12.78	202	23069	3.448	ng	97
81) Benzo(a)anthracene	13.96	228	11117	2.087	ng	# 86
83) Chrysene	14.00	228	11146	2.145	ng	# 93
88) Benzo(b)fluoranthene	15.02	252	11123m	2.775	ng	

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

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