

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102819\
 Data File : BF117567.D
 Acq On : 28 Oct 2019 17:52
 Operator : JU
 Sample : K5501-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 184-98-(0-2.2)

Manual Integrations
 APPROVED

mohammad
 10/30/2019 4:51:19 PM

Quant Time: Oct 29 17:34:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	39787	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	158961	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	79250	20.00	ng	-0.01
64) Phenanthrene-d10	11.24	188	120866	20.00	ng	0.00
76) Chrysene-d12	13.89	240	68549	20.00	ng	0.00
87) Perylene-d12	15.32	264	77311	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	204995	76.65	ng	0.00
7) Phenol-d6	6.38	99	286522	79.77	ng	0.00
23) Nitrobenzene-d5	7.29	82	174567	54.21	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	48500	70.73	ng	-0.01
45) 2-Fluorobiphenyl	9.08	172	304201	54.08	ng	-0.01
79) Terphenyl-d14	12.83	244	218660	52.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.48	163	50373	8.916	ng	98
71) Phenanthrene	11.27	178	15957	2.274	ng	99
75) Fluoranthene	12.46	202	29591	4.212	ng	98
78) Pyrene	12.69	202	25052	4.255	ng	97
81) Benzo(a)anthracene	13.88	228	10720	2.318	ng	97
83) Chrysene	13.92	228	11136	2.457	ng	97
88) Benzo(b)fluoranthene	14.92	252	13897m	3.044	ng	
90) Benzo(a)pyrene	15.27	252	8820	2.125	ng	# 92

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

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