

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100119\
 Data File : BF117286.D
 Acq On : 1 Oct 2019 19:35
 Operator : JU
 Sample : K4663-10 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-093019

Manual Integrations
 APPROVED

mohammad
 10/2/2019 3:56:44 PM

Quant Time: Oct 02 10:41:45 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	6.80	152	52032	20.00	ng	-0.02
21) Naphthalene-d8	8.09	136	206757	20.00	ng	-0.02
39) Acenaphthene-d10	9.84	164	92067	20.00	ng	-0.02
64) Phenanthrene-d10	11.33	188	117924	20.00	ng	0.00
76) Chrysene-d12	13.97	240	85265	20.00	ng	0.00
87) Perylene-d12	15.43	264	71745	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	120386	36.12	ng	-0.01
7) Phenol-d6	6.45	99	164533	38.20	ng	-0.01
23) Nitrobenzene-d5	7.37	82	97424	24.76	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	22900	29.76	ng	-0.01
45) 2-Fluorobiphenyl	9.16	172	179225	30.44	ng	-0.01
79) Terphenyl-d14	12.91	244	116225	25.48	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	33560	5.148	ng	99
75) Fluoranthene	12.54	202	32189	4.677	ng	97
78) Pyrene	12.77	202	32529	4.547	ng	99
81) Benzo(a)anthracene	13.96	228	17547	3.080	ng	# 91
83) Chrysene	14.00	228	19313	3.476	ng	# 92
88) Benzo(b)fluoranthene	15.01	252	16624m	3.825	ng	
90) Benzo(a)pyrene	15.37	252	12540	3.288	ng	# 87
92) Benzo(g,h,i)perylene	17.32	276	6810	2.249	ng	# 94

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

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