

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090419\
 Data File : BF116819.D
 Acq On : 4 Sep 2019 16:54
 Operator : HP/JU
 Sample : K4639-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DW-01

Quant Time: Sep 05 03:56:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	69503	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	277695	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	137662	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	188633	20.00	ng	0.00
76) Chrysene-d12	14.00	240	139369	20.00	ng	0.00
87) Perylene-d12	15.46	264	122866	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	458328	106.83	ng	0.00
7) Phenol-d6	6.48	99	614407	109.06	ng	0.00
23) Nitrobenzene-d5	7.40	82	360368	74.08	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	98435	106.96	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	558121	68.39	ng	0.00
79) Terphenyl-d14	12.94	244	409119	54.31	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.49	94	21571	3.458	ng	98
50) Dimethylphthalate	9.59	163	93793	9.869	ng	98
75) Fluoranthene	12.57	202	34379	3.331	ng	93
78) Pyrene	12.80	202	48480	3.913	ng	97
83) Chrysene	14.02	228	29253	3.219	ng	# 88
84) Bis(2-ethylhexyl)phthalate	13.98	149	183569	24.683	ng	99
88) Benzo(b)fluoranthene	15.03	252	33073m	4.452	ng	
90) Benzo(a)pyrene	15.40	252	13733	2.125	ng	# 1
92) Benzo(g,h,i)perylene	17.36	276	14063	2.438	ng	# 73

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

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