

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082619\
 Data File : BF116574.D
 Acq On : 26 Aug 2019 15:03
 Operator : HP/JU
 Sample : K4462-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-46-COMP

Manual Integrations
 APPROVED

Jagrut
 8/27/2019 10:30:30 AM

Quant Time: Aug 27 04:39:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 23 13:38:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	109719	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	428963	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	226364	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	380875	20.00	ng	0.00
76) Chrysene-d12	14.03	240	214161	20.00	ng	0.00
87) Perylene-d12	15.51	264	235225	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	578147	76.99	ng	0.00
7) Phenol-d6	6.50	99	780600	84.78	ng	0.00
23) Nitrobenzene-d5	7.44	82	496300	63.30	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	191553	98.55	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	854799	60.81	ng	0.00
79) Terphenyl-d14	12.98	244	745607	65.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.62	163	166368	10.534	ng	Qvalue # 99
75) Fluoranthene	12.60	202	88771	4.138	ng	96
78) Pyrene	12.83	202	108510	6.336	ng	98
81) Benzo(a)anthracene	14.02	228	45607	3.288	ng	# 86
83) Chrysene	14.06	228	45065	3.206	ng	# 89
86) Indeno(1,2,3-cd)pyrene	16.97	276	37138	3.084	ng	96
88) Benzo(b)fluoranthene	15.07	252	78140m	5.377	ng	
90) Benzo(a)pyrene	15.44	252	49486	3.853	ng	# 72
92) Benzo(g,h,i)perylene	17.42	276	41019	4.077	ng	94

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

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