

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061319\
 Data File : BF114993.D
 Acq On : 13 Jun 2019 10:26
 Operator : HP/JU
 Sample : K3159-03 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-061119-B

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:47:08 AM

Quant Time: Jun 13 18:11:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 12 15:42:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	310253	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	1177360	20.00	ng	0.00
39) Acenaphthene-d10	9.67	164	507939	20.00	ng	0.00
64) Phenanthrene-d10	11.15	188	760652	20.00	ng	-0.01
76) Chrysene-d12	13.77	240	586239	20.00	ng	-0.01
87) Perylene-d12	15.14	264	604879	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	883594	47.56	ng	0.00
7) Phenol-d6	6.29	99	1086971	48.50	ng	-0.01
23) Nitrobenzene-d5	7.20	82	631652	32.29	ng	-0.01
42) 2,4,6-Tribromophenol	10.46	330	211305	38.85	ng	-0.01
45) 2-Fluorobiphenyl	9.00	172	1112258	37.11	ng	0.00
79) Terphenyl-d14	12.73	244	775916	24.89	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.40	163	206825	5.414	ng	98
71) Phenanthrene	11.17	178	229517	5.844	ng	97
75) Fluoranthene	12.35	202	361140	8.993	ng	98
78) Pyrene	12.57	202	321921	6.028	ng	96
81) Benzo(a)anthracene	13.76	228	189674	4.324	ng	98
83) Chrysene	13.79	228	162797	3.732	ng	96
88) Benzo(b)fluoranthene	14.76	252	221215m	5.501	ng	
90) Benzo(a)pyrene	15.09	252	143598	4.139	ng	95

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

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