

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
 Data File : BF114756.D
 Acq On : 1 Jun 2019 14:52
 Operator : HP/JU
 Sample : K3085-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TANK-3

Quant Time: Jun 03 03:43:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	442280	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1645586	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	818781	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1525759	20.00	ng	0.00
76) Chrysene-d12	13.82	240	906900	20.00	ng	0.00
87) Perylene-d12	15.20	264	1043710	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1017751	40.40	ng	0.00
7) Phenol-d6	6.32	99	778190	25.64	ng	0.00
23) Nitrobenzene-d5	7.26	82	2135527	80.96	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	709533	90.95	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	3549630	80.61	ng	0.00
79) Terphenyl-d14	12.78	244	3320630	68.84	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	7.69	122	32137	2.181	ng	93
50) Dimethylphthalate	9.44	163	179992	3.155	ng	98
52) Acenaphthene	9.75	154	99188	2.100	ng	97
71) Phenanthrene	11.22	178	462843	5.947	ng	97
75) Fluoranthene	12.40	202	609768	7.566	ng	96
78) Pyrene	12.62	202	486995	6.348	ng	97

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

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