

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113493.D
 Acq On : 2 Apr 2019 6:08
 Operator : JU/SJ
 Sample : K2115-03DL 20X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW26S-GWDL

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:49:54 PM

Quant Time: Apr 02 08:34:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	99434	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	398944	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	201224	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	384013	20.00	ng	0.00
76) Chrysene-d12	13.84	240	369466	20.00	ng	0.00
87) Perylene-d12	15.26	264	268630	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	7089	1.17	ng	0.05
7) Phenol-d6	6.39	99	5945	0.77	ng	0.03
23) Nitrobenzene-d5	7.27	82	20438	3.04	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	6740	2.71	ng	0.00
45) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
79) Terphenyl-d14	12.79	244	59137	3.53	ng	0.00
Target Compounds						
27) 2,4-Dimethylphenol	7.67	122	18623m	3.492	ng	
31) Naphthalene	8.02	128	1507808	77.342	ng	99
37) 2-Methylnaphthalene	8.70	142	35516	2.920	ng	97
38) 1-Methylnaphthalene	8.80	142	53057	4.535	ng	98
52) Acenaphthene	9.77	154	24646	2.092	ng	98

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

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