

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040519\
 Data File : BF113615.D
 Acq On : 5 Apr 2019 11:21
 Operator : JU/SJ
 Sample : K2213-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S10A(2-10)COMP

Manual Integrations
 APPROVED

Sohil
 4/8/2019 9:48:11 AM

Quant Time: Apr 05 14:11:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 05 14:02:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	125054	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	465999	20.00	ng	0.00
39) Acenaphthene-d10	9.71	164	227310	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	470535	20.00	ng	-0.01
76) Chrysene-d12	13.83	240	416004	20.00	ng	0.00
87) Perylene-d12	15.23	264	329547	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	872726	119.00	ng	0.00
7) Phenol-d6	6.34	99	1132938	125.34	ng	0.00
23) Nitrobenzene-d5	7.25	82	687471	86.41	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	353808	129.57	ng	-0.01
45) 2-Fluorobiphenyl	9.04	172	1260350	89.26	ng	0.00
79) Terphenyl-d14	12.77	244	1452905	71.11	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.43	163	130982	7.817	ng	98
75) Fluoranthene	12.40	202	212794	8.493	ng	96
78) Pyrene	12.63	202	189594	5.990	ng	98
81) Benzo(a)anthracene	13.82	228	76885	3.204	ng	98
83) Chrysene	13.85	228	119905	4.929	ng	97
88) Benzo(b)fluoranthene	14.83	252	120915m	5.994	ng	

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

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