

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
 Data File : BF111942.D
 Acq On : 9 Jan 2019 11:12
 Operator : JU/SJ
 Sample : K1052-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-0103-S-DH-8A

Manual Integrations
 APPROVED

Sohil
 1/10/2019 11:45:44 AM

Quant Time: Jan 09 12:41:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	217055	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	770648	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	365078	20.00	ng	-0.01
64) Phenanthrene-d10	11.41	188	690513	20.00	ng	0.00
76) Chrysene-d12	14.05	240	495377	20.00	ng	0.00
87) Perylene-d12	15.54	264	442804	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1152708	90.86	ng	0.00
7) Phenol-d6	6.53	99	1513789	92.57	ng	0.00
23) Nitrobenzene-d5	7.45	82	965538	66.96	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	431341	90.84	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1556965	62.19	ng	0.00
79) Terphenyl-d14	12.99	244	1407803	53.95	ng	0.00
Target Compounds						
10) Phenol	6.54	94	77940	4.385	ng	82
50) Dimethylphthalate	9.63	163	182475	6.654	ng	99
78) Pyrene	12.85	202	80514	2.362	ng	97
88) Benzo(b)fluoranthene	15.10	252	60689m	2.291	ng	

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

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