

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010319\
 Data File : BF111822.D
 Acq On : 3 Jan 2019 15:49
 Operator : JU/SJ
 Sample : K1002-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-010219-D

Manual Integrations
 APPROVED

Sohil
 1/4/2019 9:56:35 AM

Quant Time: Jan 03 16:41:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 03 14:48:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	286352	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	1132369	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	582021	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	950392	20.00	ng	0.00
76) Chrysene-d12	14.06	240	629443	20.00	ng	0.00
87) Perylene-d12	15.55	264	505265	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1200830	71.48	ng	0.00
7) Phenol-d6	6.55	99	1603362	74.99	ng	0.00
23) Nitrobenzene-d5	7.46	82	1024253	52.65	ng	-0.01
42) 2,4,6-Tribromophenol	10.73	330	501913	72.98	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1809115	45.31	ng	0.00
79) Terphenyl-d14	13.01	244	1339346	40.35	ng	0.00
Target Compounds						
10) Phenol	6.55	94	75002	3.109	ng	94
50) Dimethylphthalate	9.65	163	235904	5.543	ng	97
75) Fluoranthene	12.64	202	138003	2.704	ng	95
78) Pyrene	12.87	202	135177	3.041	ng	95
88) Benzo(b)fluoranthene	15.12	252	62566m	2.119	ng	

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

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