

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100719\
 Data File : BP000550.D
 Acq On : 07 Oct 2019 18:07
 Operator : JU
 Sample : K5213-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 S-2(5-10)

Quant Time: Oct 08 06:16:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	267052	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	968629	20.00	ng	-0.01
39) Acenaphthene-d10	9.80	164	531199	20.00	ng	-0.01
64) Phenanthrene-d10	11.30	188	981158	20.00	ng	0.00
76) Chrysene-d12	13.97	240	859422	20.00	ng	0.00
87) Perylene-d12	15.45	264	973255	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1589784	95.69	ng	0.00
7) Phenol-d6	6.40	99	2114899	105.64	ng	0.00
23) Nitrobenzene-d5	7.32	82	1132548	71.41	ng	-0.01
42) 2,4,6-Tribromophenol	10.60	330	527130	93.12	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	2169320	66.08	ng	0.00
79) Terphenyl-d14	12.91	244	2535637	59.74	ng	0.00
Target Compounds						
10) Phenol	6.41	94	92012	4.489	ng	97
50) Dimethylphthalate	9.52	163	386377	10.468	ng	100
75) Fluoranthene	12.53	202	187436	3.387	ng	99
78) Pyrene	12.76	202	128880	2.174	ng	99

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

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