

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050919\
 Data File : BM020232.D
 Acq On : 10 May 2019 07:48
 Operator : JU/SJ
 Sample : K2727-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OWS-2

Quant Time: May 10 08:53:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	107512	20.00	ng	-0.02
21) Naphthalene-d8	10.57	136	378211	20.00	ng	-0.02
39) Acenaphthene-d10	14.42	164	231588	20.00	ng	-0.02
64) Phenanthrene-d10	17.17	188	555857	20.00	ng	-0.01
76) Chrysene-d12	21.35	240	596811	20.00	ng	-0.02
87) Perylene-d12	23.64	264	700714	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	967080	147.03	ng	-0.01
7) Phenol-d6	6.95	99	1170670	130.28	ng	0.00
23) Nitrobenzene-d5	8.94	82	996679	104.04	ng	-0.02
42) 2,4,6-Tribromophenol	15.91	330	651417	181.22	ng	-0.01
45) 2-Fluorobiphenyl	13.03	172	2007767	106.67	ng	-0.02
79) Terphenyl-d14	19.79	244	3078817	89.98	ng	-0.02
Target Compounds						
20) 3+4-Methylphenols	8.54	107	193724	23.408	ng	89
32) Benzoic acid	9.85	122	43634m	17.203	ng	
50) Dimethylphthalate	13.87	163	58977	2.996	ng	99

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

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