

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM073123\
 Data File : BM041204.D
 Acq On : 01 Aug 2023 01:28
 Operator : MA/JU
 Sample : 03741-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-P5A

Quant Time: Aug 01 01:59:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 31 16:29:28 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	148616	20.000	ng	0.00
21) Naphthalene-d8	10.604	136	583159	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	334995	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	703010	20.000	ng	0.00
76) Chrysene-d12	21.421	240	658453	20.000	ng	0.00
86) Perylene-d12	23.791	264	738835	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	840954	84.572	ng	0.00
7) Phenol-d6	6.975	99	1062940	82.467	ng	0.00
23) Nitrobenzene-d5	8.975	82	673445	53.704	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	407881	86.081	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	1372080	51.740	ng	0.00
79) Terphenyl-d14	19.856	244	2059869	56.562	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.927	163	123492	4.901	ng	99
84) Bis(2-ethylhexyl)phtha...	21.327	149	23977	3.390	ng	# 99

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

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