

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM073123\
 Data File : BM041207.D
 Acq On : 01 Aug 2023 03:18
 Operator : MA/JU
 Sample : 03741-13
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-P6B

Quant Time: Aug 01 03:49:42 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	181769	20.000	ng	0.00
21) Naphthalene-d8	10.604	136	698927	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	393689	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	825450	20.000	ng	0.00
76) Chrysene-d12	21.421	240	749771	20.000	ng	0.00
86) Perylene-d12	23.791	264	845122	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	778101	63.979	ng	0.00
7) Phenol-d6	6.975	99	999320	63.390	ng	0.00
23) Nitrobenzene-d5	8.969	82	620326	41.275	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	380991	68.419	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	1251231	40.149	ng	0.00
79) Terphenyl-d14	19.856	244	1792025	43.214	ng	0.00
Target Compounds						
						Qvalue
50) Dimethylphthalate	13.927	163	78856	2.663	ng	99
75) Fluoranthene	19.286	202	123180	2.391	ng	99
78) Pyrene	19.651	202	192654	3.674	ng	99
83) Chrysene	21.456	228	104126	2.118	ng	96
84) Bis(2-ethylhexyl)phtha...	21.327	149	13261	2.952	ng	97
88) Benzo(b)fluoranthene	23.068	252	105741	2.106	ng	# 92

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

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