

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM073123\
 Data File : BM041211.D
 Acq On : 01 Aug 2023 05:44
 Operator : MA/JU
 Sample : 03770-10RE
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 HOWE-2-LPRE

Quant Time: Aug 01 06:13:32 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	197761	20.000	ng	0.00
21) Naphthalene-d8	10.604	136	762142	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	427889	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	890026	20.000	ng	0.00
76) Chrysene-d12	21.421	240	800380	20.000	ng	0.00
86) Perylene-d12	23.797	264	910024	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	903513	68.283	ng	0.00
7) Phenol-d6	6.975	99	1143731	66.684	ng	0.00
23) Nitrobenzene-d5	8.969	82	708004	43.201	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	411850	68.049	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	1464833	43.246	ng	0.00
79) Terphenyl-d14	19.856	244	2133239	48.190	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.286	202	140267	2.525	ng	98
78) Pyrene	19.651	202	136058	2.431	ng	98
84) Bis(2-ethylhexyl)phtha...	21.333	149	16569	3.023	ng	# 84

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM073123\
Data File : BM041211.D
Acq On : 01 Aug 2023 05:44
Operator : MA/JU
Sample : 03770-10RE
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
HOWE-2-LPRE

Quant Time: Aug 01 06:13:32 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

