

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080823\
 Data File : BM041338.D
 Acq On : 08 Aug 2023 21:05
 Operator : MA/JU
 Sample : 03929-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-PP19

Quant Time: Aug 09 01:20:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 17:25:06 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	131755	20.000	ng	0.00
21) Naphthalene-d8	10.580	136	520369	20.000	ng	-0.01
39) Acenaphthene-d10	14.445	164	298868	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	620776	20.000	ng	-0.01
76) Chrysene-d12	21.415	240	587802	20.000	ng	0.00
86) Perylene-d12	23.780	264	676114	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.351	112	657917	74.632	ng	0.00
7) Phenol-d6	6.963	99	834149	72.998	ng	0.00
23) Nitrobenzene-d5	8.957	82	541376	48.382	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	322718	76.341	ng	0.00
45) 2-Fluorobiphenyl	13.063	172	1055660	44.620	ng	0.00
79) Terphenyl-d14	19.845	244	1443210	44.393	ng	0.00
Target Compounds						
78) Pyrene	19.639	202	129598	3.153	ng	99
83) Chrysene	21.450	228	83259	2.160	ng	98
84) Bis(2-ethylhexyl)phtha...	21.321	149	9838	2.930	ng	# 98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080823\
Data File : BM041338.D
Acq On : 08 Aug 2023 21:05
Operator : MA/JU
Sample : 03929-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-PP19

Quant Time: Aug 09 01:20:40 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 07 17:25:06 2023
Response via : Initial Calibration

