

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032967.D
 Acq On : 10 Nov 2021 23:58
 Operator : CG/JU
 Sample : M4492-01
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG1R8

Quant Time: Nov 11 09:29:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 13:31:56 2021
 Response via : Initial Calibration

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

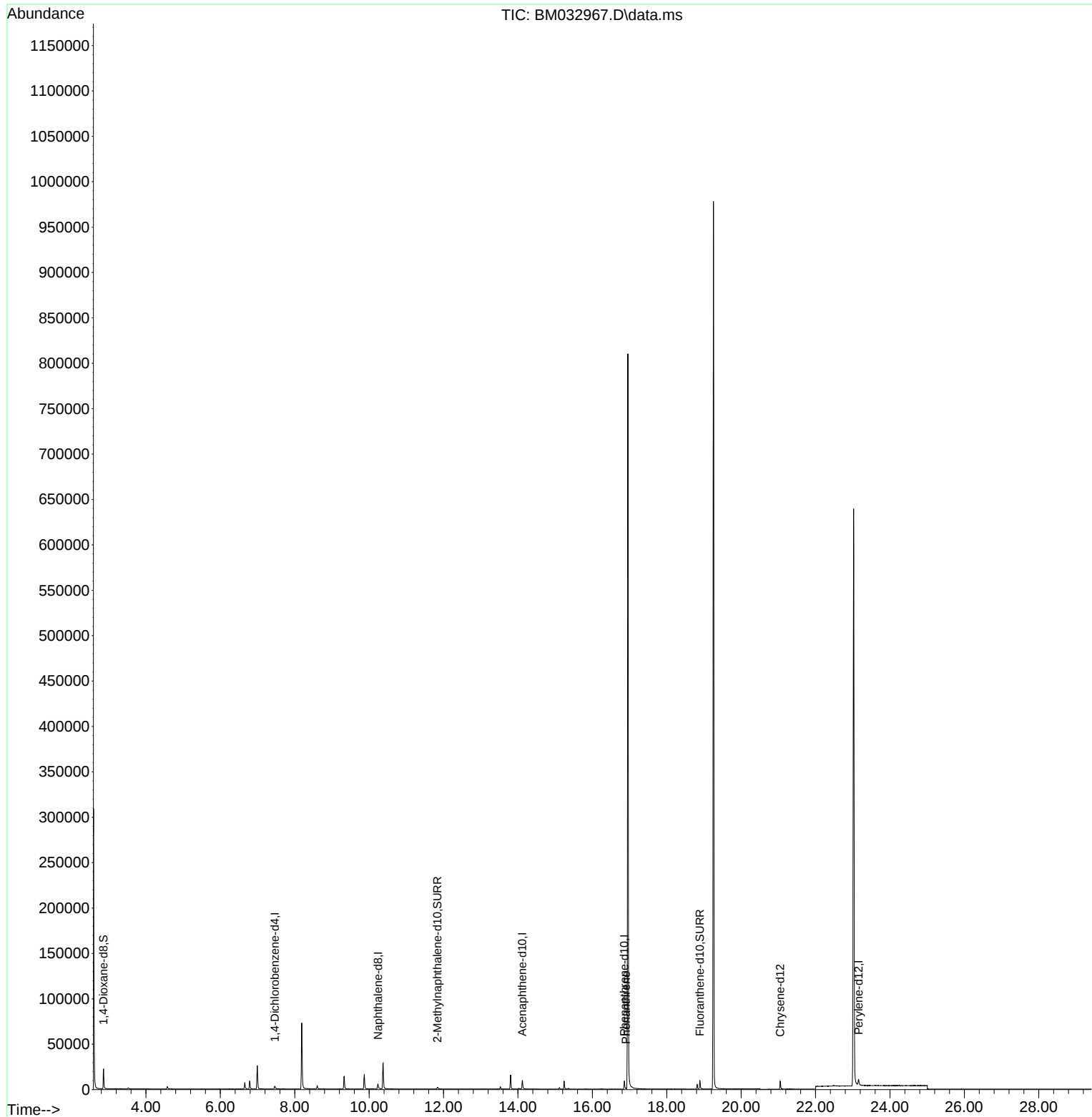
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.462	152	2362	0.400	ng/ul	0.00
4) Naphthalene-d8	10.235	136	8666	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.117	164	5235	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	16.862	188	10919	0.400	ng/ul	0.00
17) Chrysene-d12	21.049	240	8814	0.400	ng/ul	0.00
23) Perylene-d12	23.154	264	8498	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.860	96	13030	4.320	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.836	152	3550	0.288	ng/ul	0.00
18) Fluoranthene-d10	18.888	212	10943	0.410	ng/ul	0.00
Target Compounds						
15) Phenanthrene	16.900	178	807	0.022	ng/ul#	Qvalue 83

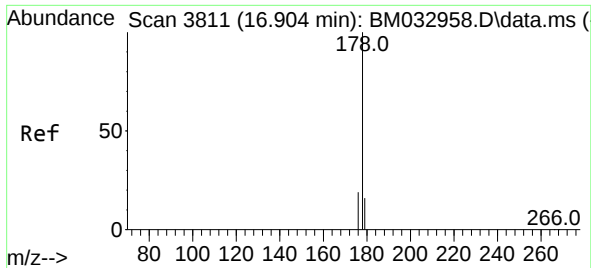
(#) = qualifier out of range (m) = manual integration (+) = signals summed

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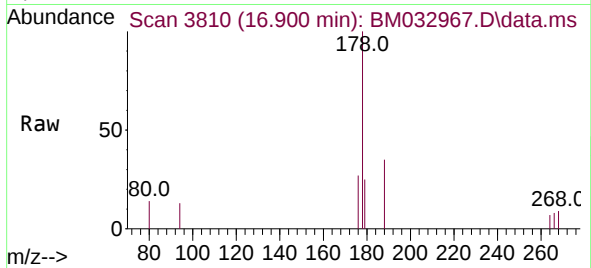
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#15
 Phenanthrene
 Concen: 0.022 ng/ul
 RT: 16.900 min Scan# 3810
 Delta R.T. -0.008 min
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Tgt Ion:	178	Resp:	807
Ion Ratio	100	Lower	Upper
178	100		
179	24.9	12.9	19.3#
176	26.7	16.0	24.0#

