

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092021\
 Data File : BN016363.D
 Acq On : 21 Sep 2021 03:46
 Operator : CG/JU
 Sample : M3729-10
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-TT-VPB-RW7-GW-790-792

Manual Integrations
 APPROVED

Quant Time: Sep 21 04:41:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 20 11:09:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.679	152	25295	0.400	ng	0.00
7) Naphthalene-d8	10.445	136	113786	0.400	ng	#-0.01
13) Acenaphthene-d10	14.307	164	63398	0.400	ng	0.00
19) Phenanthrene-d10	17.054	188	129536	0.400	ng	0.00
29) Chrysene-d12	21.259	240	139378	0.400	ng	# 0.00
35) Perylene-d12	23.515	264	142673	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.319	112	3601	0.057	ng	0.00
5) Phenol-d6	6.868	99	2780	0.036	ng	0.00
8) Nitrobenzene-d5	8.823	82	16796	0.248	ng	0.00
11) 2-Methylnaphthalene-d10	12.043	152	40340	0.225	ng	0.00
14) 2,4,6-Tribromophenol	15.804	330	3246	0.174	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	57475	0.232	ng	-0.01
27) Fluoranthene-d10	19.093	212	89089	0.238	ng	0.00
31) Terphenyl-d14	19.703	244	91513	0.283	ng	0.00
Target Compounds						Qvalue
18) Fluorene	15.360	166	6436	0.028	ng	# 88
24) Pentachlorophenol	16.713	266	527	0.022	ng	94
25) Phenanthrene	17.091	178	97035m	0.265	ng	
26) Anthracene	17.188	178	9112	0.026	ng	# 82
28) Fluoranthene	19.123	202	21735	0.052	ng	97
30) Pyrene	19.485	202	30947	0.070	ng	99
32) Benzo(a)anthracene	21.238	228	13158	0.029	ng	# 78
33) Chrysene	21.291	228	14059	0.032	ng	# 82
34) Bis(2-ethylhexyl)phtha...	21.196	149	349709	1.448	ng	99
36) Indeno(1,2,3-cd)pyrene	25.809	276	13614	0.030	ng	# 95
37) Benzo(b)fluoranthene	22.834	252	23818	0.053	ng	# 12
38) Benzo(k)fluoranthene	22.879	252	14381	0.033	ng	# 1
39) Benzo(a)pyrene	23.416	252	11824	0.029	ng	# 1
40) Dibenzo(a,h)anthracene	25.829	278	10172	0.026	ng	# 38
41) Benzo(g,h,i)perylene	26.510	276	11718	0.030	ng	# 79

mohammad

9/21/2021 11:30:44 AM

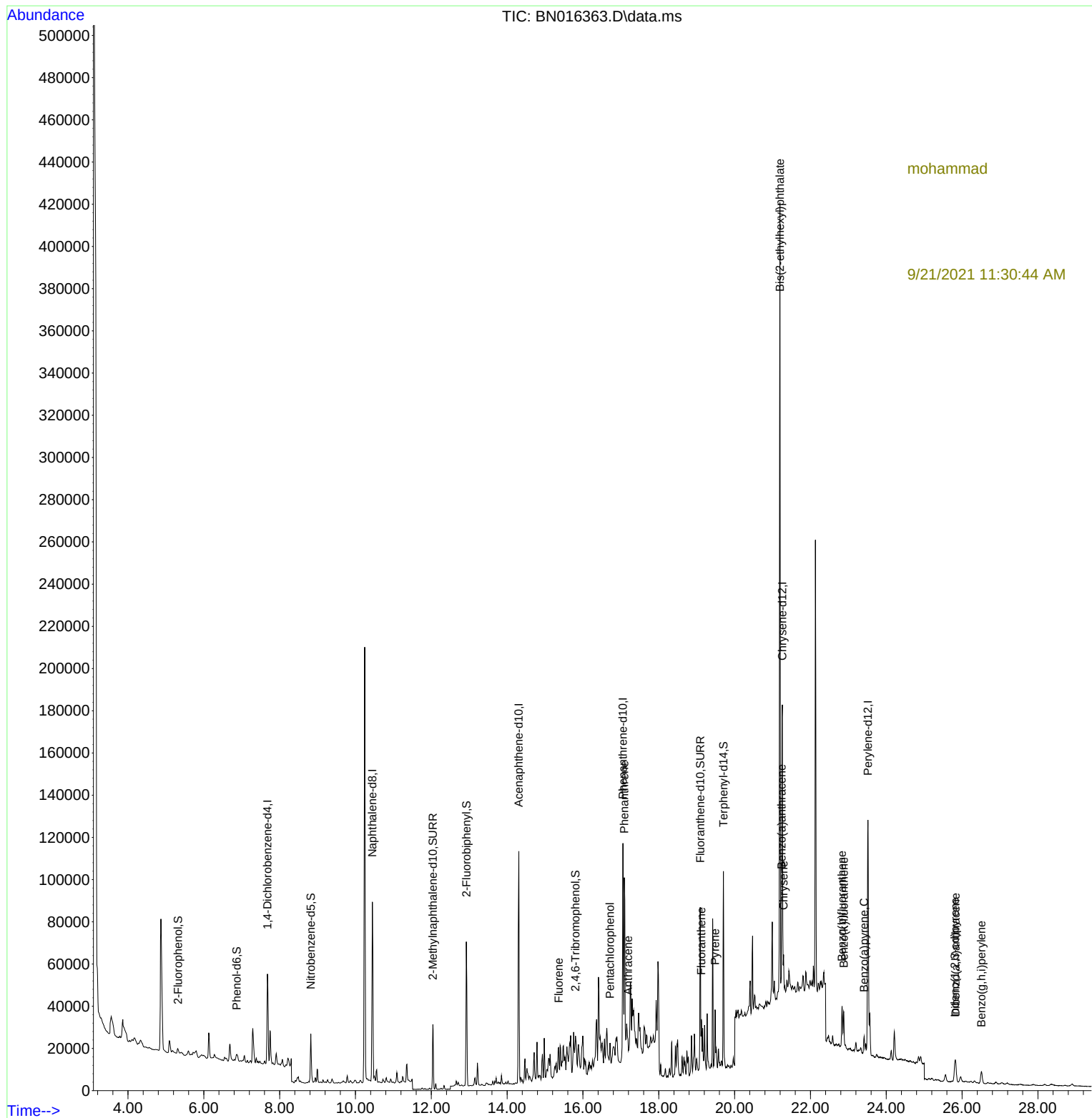
(#) = qualifier out of range (m) = manual integration (+) = signals summed

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