

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE103124\
 Data File : BE101440.D
 Acq On : 31 Oct 2024 19:24
 Operator : RC/JU
 Sample : P4611-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/04/2024

Quant Time: Nov 01 06:16:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE102824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 29 01:26:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.573	152	129070	20.000	ng	0.00
21) Naphthalene-d8	10.346	136	590260	20.000	ng	0.00
39) Acenaphthene-d10	14.182	164	401080	20.000	ng	0.00
64) Phenanthrene-d10	16.920	188	855474	20.000	ng	0.00
76) Chrysene-d12	21.080	240	882545	20.000	ng	0.00
86) Perylene-d12	23.378	264	1145405	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.322	112	722080	98.053	ng	0.00
7) Phenol-d6	6.956	99	991118	97.062	ng	0.00
23) Nitrobenzene-d5	8.783	82	591310	62.921	ng	0.00
42) 2,4,6-Tribromophenol	15.693	330	699114	99.452	ng	0.00
45) 2-Fluorobiphenyl	12.808	172	1615692	68.222	ng	0.00
79) Terphenyl-d14	19.517	244	2935980	76.566	ng	0.00
Target Compounds						Qvalue
78) Pyrene	19.335	202	97831	2.016	ng	98
83) Chrysene	21.098	228	133678	2.787	ng	98
88) Benzo(b)fluoranthene	22.661	252	285445m	4.782	ng	
90) Benzo(a)pyrene	23.266	252	147436	2.824	ng	# 87

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

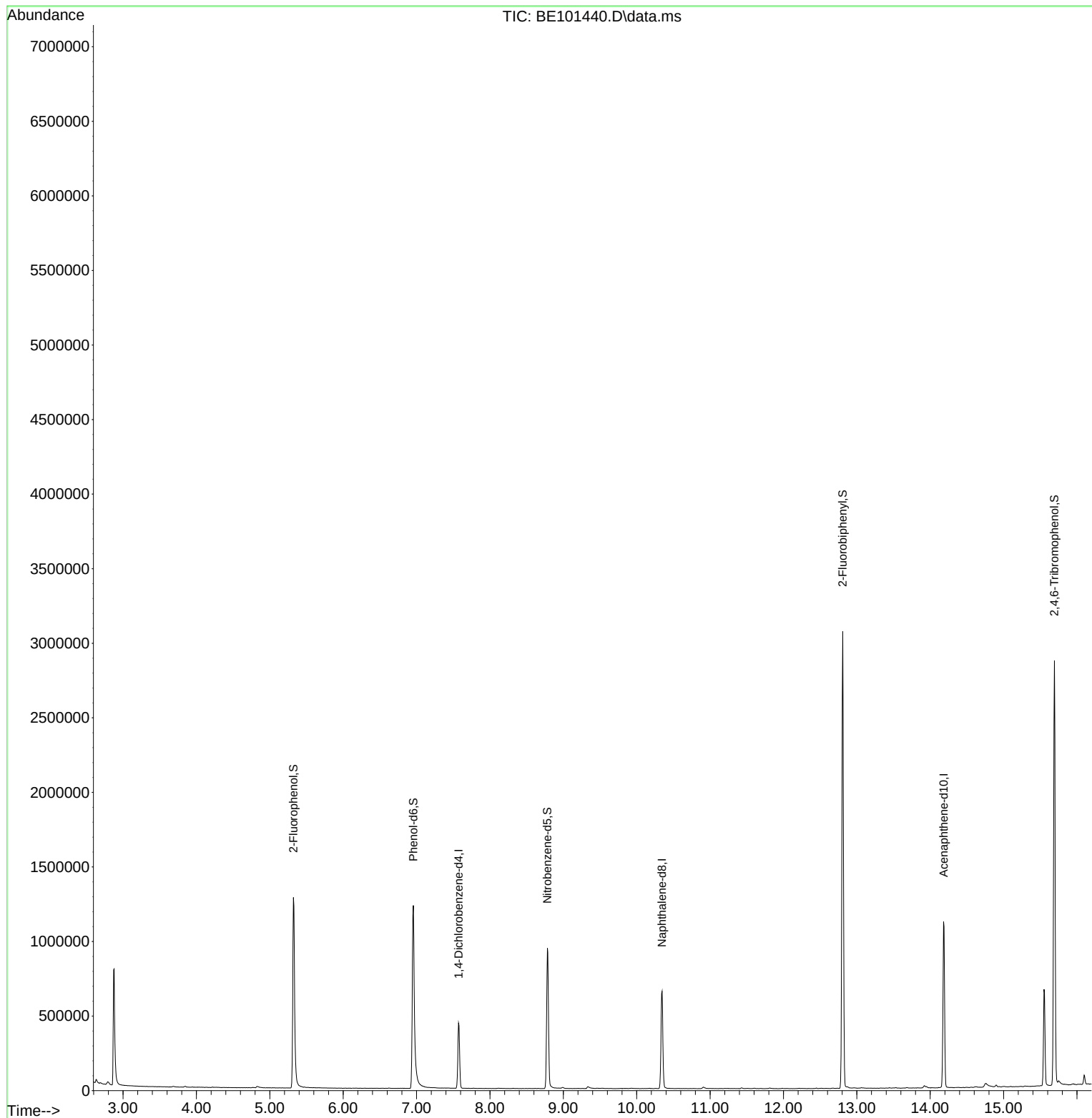
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