

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092520\
 Data File : BP003391.D
 Acq On : 25 Sep 2020 20:41
 Operator : CG/JU
 Sample : L4127-01DL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B-3(9-11)DL

Manual Integrations
 APPROVED

mohammad
 9/28/2020 11:11:23 AM

Quant Time: Sep 26 01:25:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 12:48:08 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	161093	20.00	ng	0.00
21) Naphthalene-d8	10.352	136	697358	20.00	ng	# 0.00
39) Acenaphthene-d10	14.222	164	425451	20.00	ng	0.00
64) Phenanthrene-d10	16.981	188	818043	20.00	ng	0.00
76) Chrysene-d12	21.080	240	649455	20.00	ng	0.01
86) Perylene-d12	23.280	264	679357	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	146760	15.91	ng	0.00
7) Phenol-d6	6.781	99	185582	15.09	ng	0.01
23) Nitrobenzene-d5	8.722	82	107755	9.08	ng	0.00
42) 2,4,6-Tribromophenol	15.728	330	70943	10.93	ng	0.00
45) 2-Fluorobiphenyl	12.840	172	319341	10.98	ng	0.00
79) Terphenyl-d14	19.557	244	363859	11.67	ng	0.00
Target Compounds						Qvalue
58) Fluorene	15.281	166	83089	2.70	ng	99
71) Phenanthrene	17.022	178	977462	21.96	ng	100
72) Anthracene	17.110	178	287991	6.57	ng	99
75) Fluoranthene	19.004	202	1379880	24.66	ng	98
78) Pyrene	19.357	202	1189276	29.36	ng	100
81) Benzo(a)anthracene	21.063	228	609284	14.51	ng	99
83) Chrysene	21.116	228	531062	13.17	ng	97
87) Indeno(1,2,3-cd)pyrene	25.521	276	339204	6.59	ng	95
88) Benzo(b)fluoranthene	22.627	252	691941	16.13	ng	97
89) Benzo(k)fluoranthene	22.663	252	214594m	5.22	ng	
90) Benzo(a)pyrene	23.186	252	506738	12.53	ng	95
91) Dibenzo(a,h)anthracene	25.521	278	92621	2.18	ng	# 84
92) Benzo(g,h,i)perylene	26.204	276	324610	7.65	ng	97

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

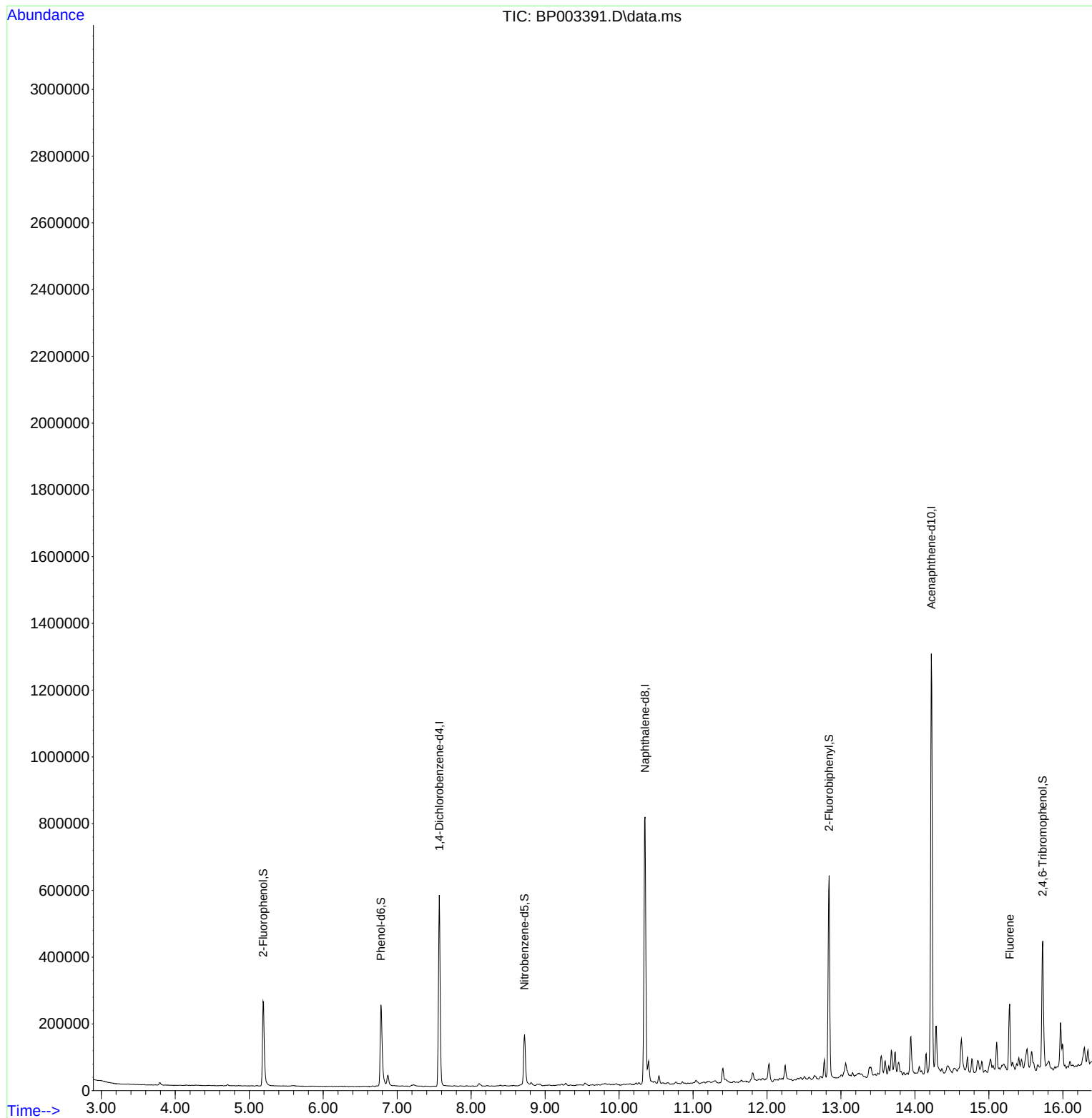
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