

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP093020\
 Data File : BP003431.D
 Acq On : 01 Oct 2020 01:09
 Operator : CG/JU
 Sample : L4193-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B-12(9-11)

Manual Integrations
 APPROVED

mohammad
 10/1/2020 1:34:07 PM

Quant Time: Oct 01 02:20:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 30 17:14:02 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.540	152	98368	20.00	ng	0.00
21) Naphthalene-d8	10.316	136	410295	20.00	ng	# 0.00
39) Acenaphthene-d10	14.198	164	239932	20.00	ng	0.00
64) Phenanthrene-d10	16.963	188	526257	20.00	ng	# 0.02
76) Chrysene-d12	21.163	240	523331	20.00	ng	0.10
86) Perylene-d12	23.680	264	550601m	20.00	ng	0.42
System Monitoring Compounds						
5) 2-Fluorophenol	5.158	112	521530	92.58	ng	0.00
7) Phenol-d6	6.746	99	696031	92.69	ng	0.00
23) Nitrobenzene-d5	8.693	82	460512	65.94	ng	0.00
42) 2,4,6-Tribromophenol	15.710	330	262599	71.76	ng	0.02
45) 2-Fluorobiphenyl	12.810	172	975357	59.45	ng	0.00
79) Terphenyl-d14	19.569	244	1386080	55.18	ng	0.04
Target Compounds						Qvalue
31) Naphthalene	10.369	128	94170	4.34	ng	98
49) Acenaphthylene	13.916	152	149309	6.59	ng	94
50) Dimethylphthalate	13.657	163	79493	4.15	ng	# 96
52) Acenaphthene	14.263	154	68795	4.99	ng	98
55) Dibenzofuran	14.604	168	75949	3.45	ng	# 88
58) Fluorene	15.257	166	98023	5.64	ng	# 92
71) Phenanthrene	17.010	178	1142111	39.89	ng	99
72) Anthracene	17.098	178	402015	14.25	ng	98
73) Carbazole	17.381	167	147663	5.51	ng	# 91
75) Fluoranthene	19.022	202	2247203	62.42	ng	99
78) Pyrene	19.381	202	2016497	61.78	ng	100
81) Benzo(a)anthracene	21.163	228	1251494	36.99	ng	# 92
83) Chrysene	21.192	228	980274m	30.16	ng	
88) Benzo(b)fluoranthene	22.880	252	2181745m	62.75	ng	
89) Benzo(k)fluoranthene	23.521	252	2083648m	62.53	ng	
90) Benzo(a)pyrene	23.521	252	2106470m	64.25	ng	
92) Benzo(g,h,i)perylene	26.463	276	941764m	27.39	ng	

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

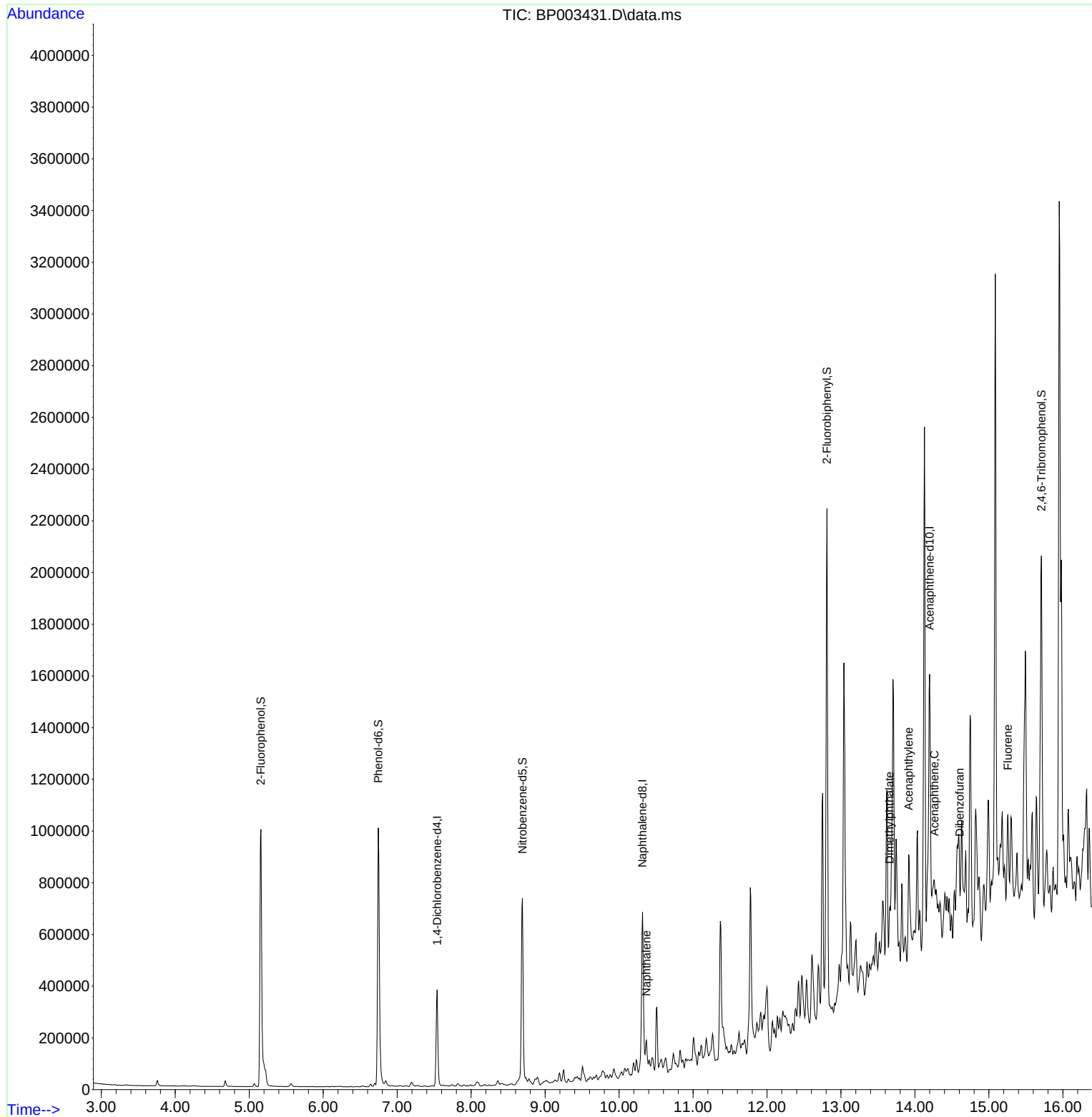
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