

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110524\
 Data File : BE101476.D
 Acq On : 5 Nov 2024 13:20
 Operator : RC/JU
 Sample : P4675-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 COMP-2

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/06/2024
 Supervised By :mohammad ahmed 11/06/2024

Quant Time: Nov 05 13:06:20 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	7.570	152	54196	20.000	ng	0.00
21) Naphthalene-d8	10.337	136	232476	20.000	ng	0.00
39) Acenaphthene-d10	14.180	164	158291	20.000	ng	0.00
64) Phenanthrene-d10	16.918	188	377849	20.000	ng	0.00
76) Chrysene-d12	21.077	240	468893	20.000	ng	0.00
86) Perylene-d12	23.363	264	597551	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.319	112	281954	91.182	ng	0.00
7) Phenol-d6	6.947	99	372313	86.834	ng	0.00
23) Nitrobenzene-d5	8.774	82	212712	57.469	ng	-0.01
42) 2,4,6-Tribromophenol	15.684	330	250898	90.436	ng	0.00
45) 2-Fluorobiphenyl	12.799	172	578229	61.864	ng	0.00
79) Terphenyl-d14	19.509	244	1353479	66.435	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	16.953	178	64575	3.569	ng	98
75) Fluoranthene	18.962	202	134883	6.139	ng	95
78) Pyrene	19.327	202	116413	4.515	ng	96
81) Benzo(a)anthracene	21.054	228	74209	2.816	ng	95
83) Chrysene	21.107	228	68005	2.669	ng	95
88) Benzo(b)fluoranthene	22.652	252	72955m	2.343	ng	
90) Benzo(a)pyrene	23.251	252	68692	2.522	ng	99

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

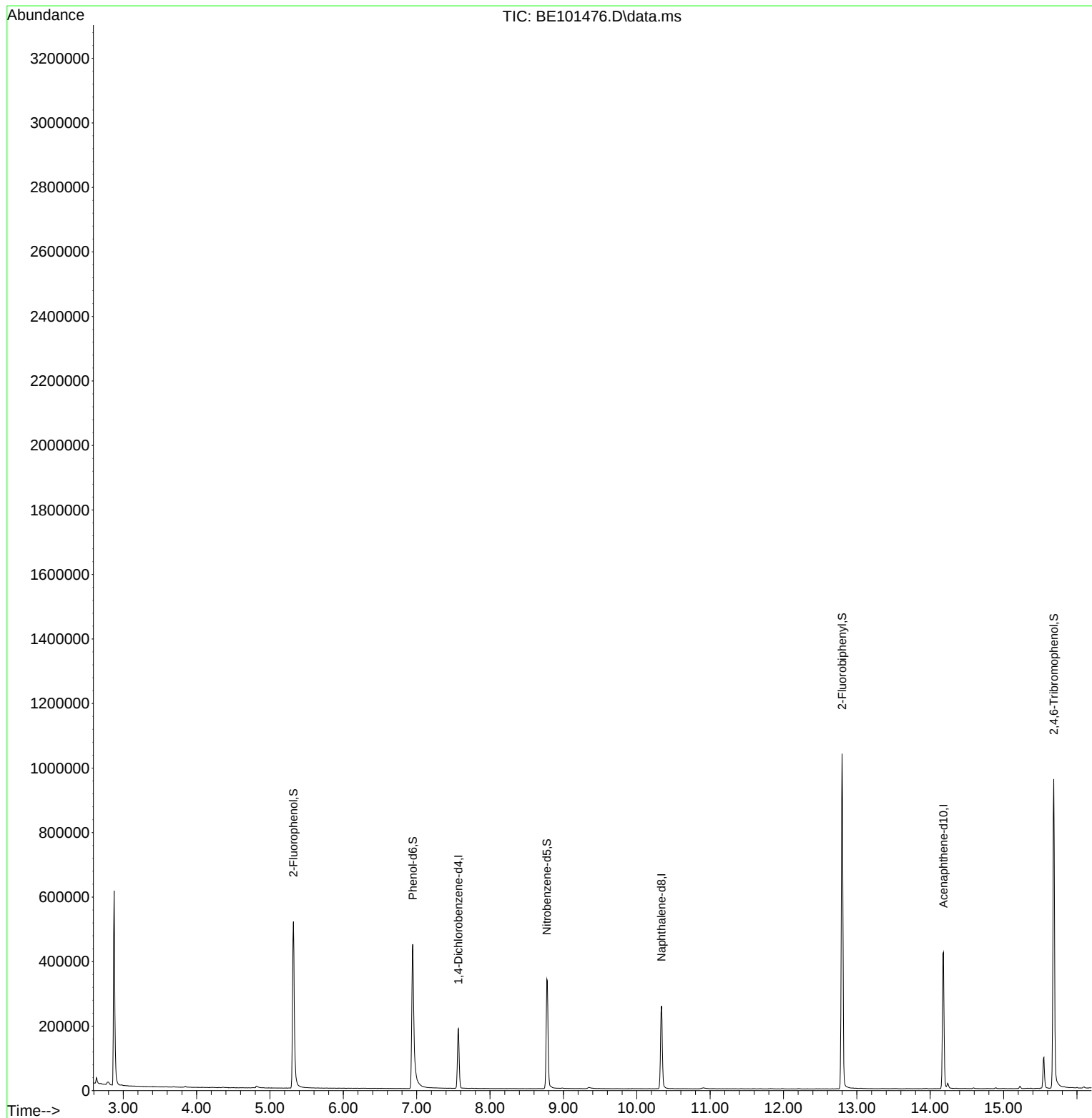
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