

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110724\
 Data File : BE101534.D
 Acq On : 7 Nov 2024 16:54
 Operator : RC/JU
 Sample : P4711-06DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 CF-613-COMP-17DL

Quant Time: Nov 07 16:58:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 00:01:20 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 11/08/2024
 Supervised By :mohammad ahmed 11/08/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.559	152	65686	20.000	ng	0.00
21) Naphthalene-d8	10.326	136	304849	20.000	ng	0.00
39) Acenaphthene-d10	14.169	164	203453	20.000	ng	0.00
64) Phenanthrene-d10	16.907	188	447469	20.000	ng	0.00
76) Chrysene-d12	21.072	240	473017	20.000	ng	0.00
86) Perylene-d12	23.358	264	620522	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.314	112	50699	13.647	ng	0.00
7) Phenol-d6	6.942	99	66968	13.148	ng	0.00
23) Nitrobenzene-d5	8.769	82	39264	8.136	ng	0.00
42) 2,4,6-Tribromophenol	15.679	330	47113	12.306	ng	0.00
45) 2-Fluorobiphenyl	12.788	172	117782	9.453	ng	-0.01
79) Terphenyl-d14	19.498	244	235065	11.000	ng	0.00
Target Compounds						Qvalue
49) Acenaphthylene	13.904	152	47345	2.870	ng	99
52) Acenaphthene	14.233	154	45636	4.336	ng	98
55) Dibenzofuran	14.574	168	46388	2.737	ng	98
58) Fluorene	15.214	166	82369	5.823	ng	95
71) Phenanthrene	16.953	178	1243943	57.157	ng	100
72) Anthracene	17.036	178	350066	16.288	ng	98
73) Carbazole	17.365	167	44354	2.058	ng	98
75) Fluoranthene	18.963	202	2080746	74.568	ng	98
78) Pyrene	19.327	202	1787674	66.422	ng	98
81) Benzo(a)anthracene	21.055	228	1125386	40.060	ng	100
83) Chrysene	21.107	228	955066	35.767	ng	98
87) Indeno(1,2,3-cd)pyrene	25.679	276	751130	17.615	ng	97
88) Benzo(b)fluoranthene	22.653	252	1245683	36.590	ng	98
89) Benzo(k)fluoranthene	22.688	252	657408m	21.272	ng	
90) Benzo(a)pyrene	23.252	252	1037703	35.306	ng	99
91) Dibenzo(a,h)anthracene	25.661	278	190450m	5.363	ng	
92) Benzo(g,h,i)perylene	26.431	276	702021	19.457	ng	97

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

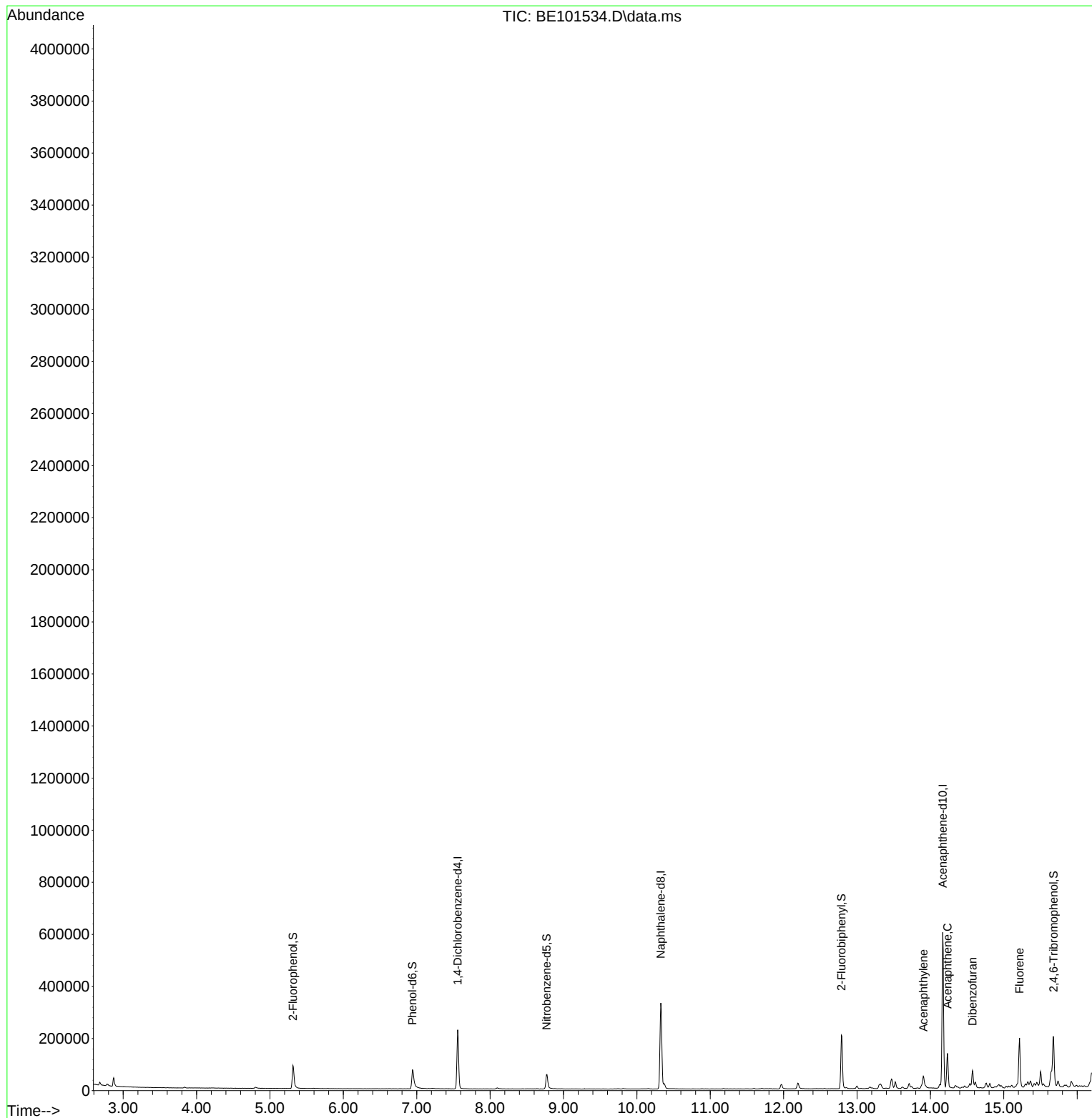
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