

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042121\
 Data File : BG048813.D
 Acq On : 21 Apr 2021 9:42
 Operator : CG/JU
 Sample : PB135874BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB135874BL

Quant Time: Apr 21 11:19:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG041421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 20 17:33:30 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	18289	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	71967	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	51811	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	131604	20.00	ng	0.00
76) Chrysene-d12	21.77	240	142536	20.00	ng	0.00
86) Perylene-d12	25.10	264	139506	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.61	112	154887	141.64	ng	0.00
7) Phenol-d6	7.23	99	192551	124.49	ng	0.00
23) Nitrobenzene-d5	9.23	82	144652	89.99	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	88335	127.16	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	324936	90.57	ng	0.00
79) Terphenyl-d14	20.09	244	703855	84.99	ng	0.00

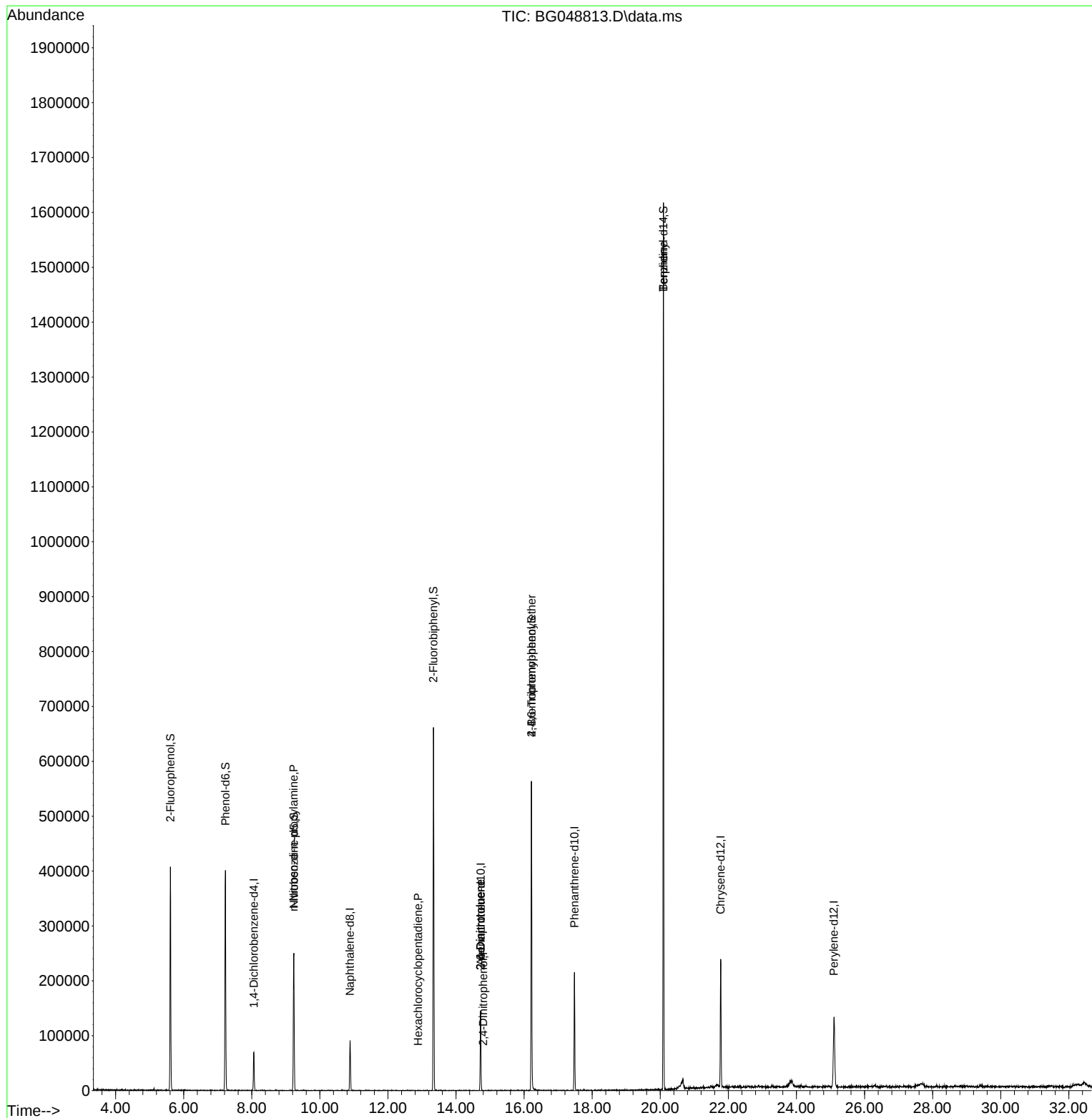
Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
19) n-Nitroso-di-n-propylamine	9.23	70	22829	21.877	ng	# 80
41) Hexachlorocyclopentadiene	12.88	237	59	5.919	ng	# 31
51) 2,6-Dinitrotoluene	14.72	165	6814	7.600	ng	# 10
54) 2,4-Dinitrophenol	14.80	184	57	5.380	ng	# 8
57) 2,4-Dinitrotoluene	14.72	165	6814	5.233	ng	# 15
67) 4-Bromophenyl-phenylether	16.21	248	7027	4.639	ng	# 1
77) Benzidine	20.09	184	14133	3.651	ng	# 90

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

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#1
1,4-Dichlorobenzene-d4
Concen: 20.000 ng

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