

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011921\
 Data File : BF123015.D
 Acq On : 19 Jan 2021 12:50
 Operator : JU/CG
 Sample : M1141-23
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Jan 19 13:17:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF011621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 18 13:14:08 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	186169	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	767445	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	365362	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	663859	20.00	ng	0.00
76) Chrysene-d12	14.08	240	660819	20.00	ng	0.00
86) Perylene-d12	15.57	264	646751	20.00	ng	0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.53	112	1464798	118.93	ng	0.02
7) Phenol-d6	6.52	99	1835768	110.91	ng	0.00
23) Nitrobenzene-d5	7.46	82	1313803	88.09	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	621382	150.39	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	2075256	79.54	ng	0.00
79) Terphenyl-d14	13.03	244	3776293	104.78	ng	0.02

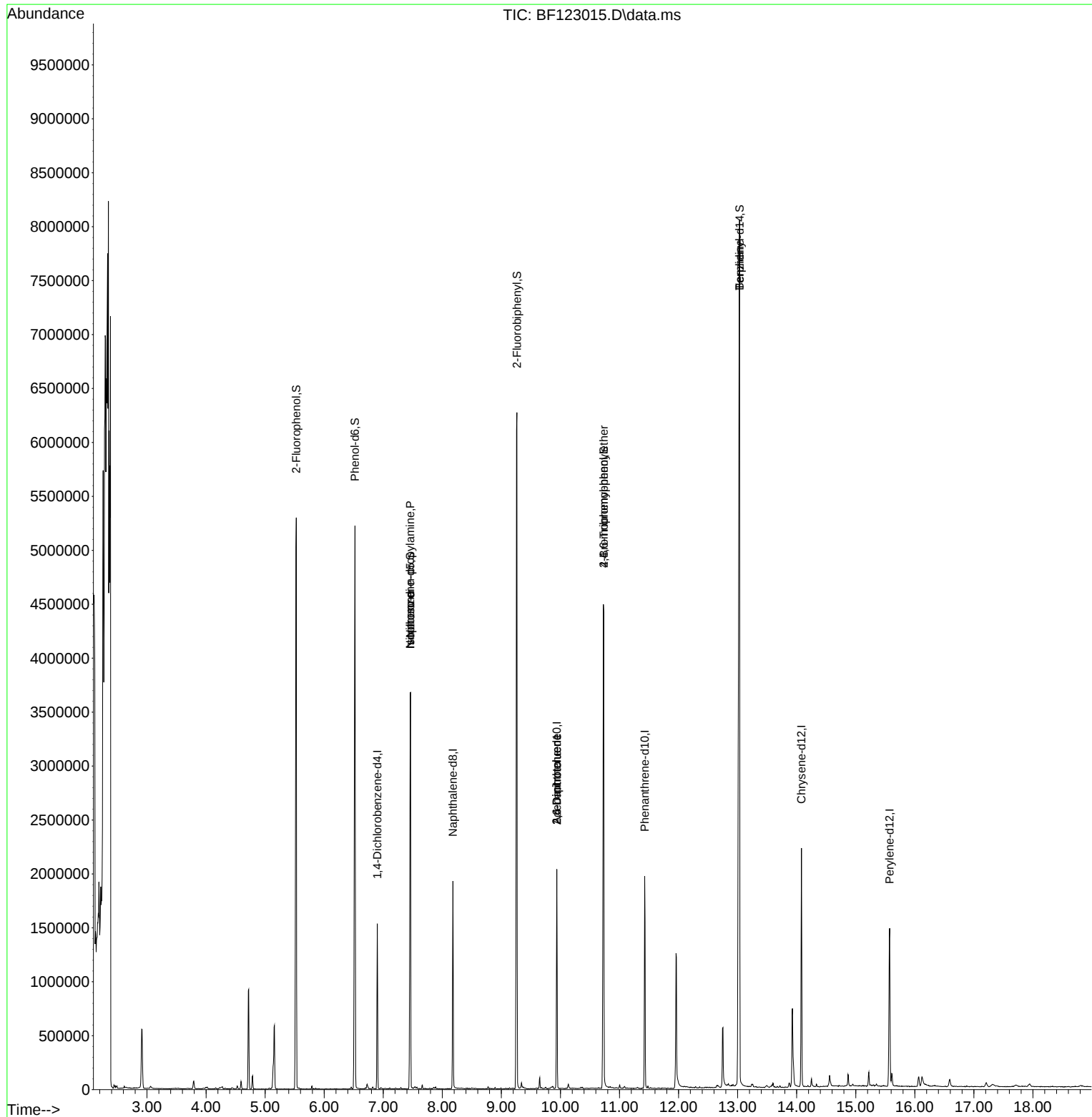
Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Benzaldehyde	6.45	77	401	Below Cal	#	81
19) n-Nitroso-di-n-propylamine	7.46	70	178507	18.205	ng	# 83
25) Isophorone	7.46	82	1313788	49.375	ng	# 65
51) 2,6-Dinitrotoluene	9.94	165	47009	7.579	ng	# 31
57) 2,4-Dinitrotoluene	9.94	165	47009	5.892	ng	# 25
67) 4-Bromophenyl-phenylether	10.73	248	38179	4.704	ng	# 1
77) Benzidine	13.03	184	52503	3.000	ng	# 89

(#) = 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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