

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070720\
 Data File : BF120784.D
 Acq On : 7 Jul 2020 13:42
 Operator : JU/CG
 Sample : L3215-16
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2

Quant Time: Jul 07 16:14:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 02 18:51:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	154186	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	590927	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	306147	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	547815	20.00	ng	0.00
76) Chrysene-d12	14.00	240	420757	20.00	ng	0.00
86) Perylene-d12	15.47	264	347018	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	188303	18.85	ng	0.00
7) Phenol-d6	6.47	99	399297	31.30	ng	0.00
23) Nitrobenzene-d5	7.42	82	951724	93.31	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	477534	155.26	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	1769891	87.47	ng	0.00
79) Terphenyl-d14	12.96	244	2157273	100.38	ng	0.00
Target Compounds						
10) Phenol	6.49	94	29206	2.226	ng	88
27) 2,4-Dimethylphenol	7.79	122	353479	40.257	ng	99
31) Naphthalene	8.16	128	2070143	65.551	ng	97
37) 2-Methylnaphthalene	8.85	142	93366	4.548	ng	# 80
38) 1-Methylnaphthalene	8.95	142	323041	16.799	ng	# 98
50) Dimethylphthalate	9.60	163	67080	3.004	ng	# 91
73) Carbazole	11.60	167	102299	3.544	ng	99

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

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