

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082420\
 Data File : BM027202.D
 Acq On : 24 Aug 2020 20:02
 Operator : JU/CG
 Sample : L3734-07RE 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EROC-BF-1FRE

Quant Time: Aug 25 01:33:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082120.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 15:18:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	216509	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	809337	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	513922	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1063241	20.00	ng	0.00
76) Chrysene-d12	21.19	240	999216	20.00	ng	0.00
86) Perylene-d12	23.37	264	1104673	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	101238	9.12	ng	0.00
7) Phenol-d6	6.79	99	464429	30.46	ng	0.00
23) Nitrobenzene-d5	8.75	82	411993	21.00	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	21065	2.53	ng	0.00
45) 2-Fluorobiphenyl	12.86	172	842438	21.53	ng	0.00
79) Terphenyl-d14	19.63	244	1055126	19.73	ng	0.00
Target Compounds						
71) Phenanthrene	17.03	178	182929	3.095	ng	99
75) Fluoranthene	19.05	202	279332	3.783	ng	97
78) Pyrene	19.42	202	266232	4.379	ng	99
81) Benzo(a)anthracene	21.17	228	139393	2.153	ng	# 90
83) Chrysene	21.22	228	138892	2.209	ng	# 92
84) Bis(2-ethylhexyl)phthalate	21.13	149	438423	13.861	ng	99
88) Benzo(b)fluoranthene	22.72	252	177325	2.407	ng	# 79

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

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