

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082420\
 Data File : BM027205.D
 Acq On : 24 Aug 2020 21:51
 Operator : JU/CG
 Sample : L3734-02RE 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EROC-BF-1ARE

Manual Integrations
 APPROVED

mohammad
 8/25/2020 4:18:31 PM

Quant Time: Aug 25 02:02:09 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	209701	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	791158	20.00	ng	0.00
39) Acenaphthene-d10	14.24	164	489985	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1008600	20.00	ng	0.00
76) Chrysene-d12	21.19	240	954252	20.00	ng	0.00
86) Perylene-d12	23.38	264	1060548	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	189571	17.62	ng	0.00
7) Phenol-d6	6.79	99	604949	40.96	ng	0.00
23) Nitrobenzene-d5	8.75	82	531529	27.72	ng	0.00
42) 2,4,6-Tribromophenol	15.74	330	55073	6.94	ng	0.00
45) 2-Fluorobiphenyl	12.86	172	1091342	29.26	ng	0.00
79) Terphenyl-d14	19.63	244	1357847	26.58	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.71	163	92251	2.319	ng	100
71) Phenanthrene	17.03	178	183659	3.276	ng	98
75) Fluoranthene	19.06	202	361439	5.161	ng	100
78) Pyrene	19.42	202	325771	5.610	ng	98
81) Benzo(a)anthracene	21.17	228	174757	2.826	ng	# 89
83) Chrysene	21.23	228	177285	2.952	ng	# 93
84) Bis(2-ethylhexyl)phthalate	21.13	149	238609	7.899	ng	97
88) Benzo(b)fluoranthene	22.73	252	250280m	3.539	ng	
90) Benzo(a)pyrene	23.28	252	167752	2.598	ng	# 79

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

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