

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012920\
 Data File : BM024656.D
 Acq On : 29 Jan 2020 18:41
 Operator : CG/JU
 Sample : L1013-07 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PL-05-012820

Manual Integrations
 APPROVED

mohammad
 1/30/2020 10:19:09 AM

Quant Time: Jan 30 04:02:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 23 14:57:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	440559	20.00	ng	0.00
21) Naphthalene-d8	10.20	136	1683333	20.00	ng	0.00
39) Acenaphthene-d10	14.09	164	1018027	20.00	ng	0.00
64) Phenanthrene-d10	16.84	188	1802749	20.00	ng	0.00
76) Chrysene-d12	21.06	240	1419159	20.00	ng	0.00
87) Perylene-d12	23.17	264	1679972	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.09	112	1017548	43.92	ng	0.00
7) Phenol-d6	6.65	99	1314364	41.19	ng	0.00
23) Nitrobenzene-d5	8.58	82	842632	24.55	ng	0.00
42) 2,4,6-Tribromophenol	15.59	330	614355	37.10	ng	0.00
45) 2-Fluorobiphenyl	12.70	172	2218397	29.63	ng	0.00
79) Terphenyl-d14	19.50	244	2307481	30.35	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.56	163	366395	4.451	ng	98
71) Phenanthrene	16.89	178	330332	3.384	ng	97
75) Fluoranthene	18.92	202	806282	6.617	ng	98
78) Pyrene	19.28	202	750519	8.022	ng	99
81) Benzo(a)anthracene	21.04	228	398040	4.041	ng	# 90
83) Chrysene	21.09	228	376606m	4.007	ng	
86) Indeno(1,2,3-cd)pyrene	25.25	276	294659	2.876	ng	98
88) Benzo(b)fluoranthene	22.56	252	567336m	5.207	ng	
90) Benzo(a)pyrene	23.09	252	422177	4.262	ng	# 85
92) Benzo(g,h,i)perylene	25.88	276	302199	3.540	ng	95

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

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