

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060420\
 Data File : BF120532.D
 Acq On : 4 Jun 2020 14:39
 Operator : JU/CG
 Sample : L2839-16
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM111-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/5/2020 11:57:16 AM

Quant Time: Jun 04 16:46:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 04 16:38:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	175810	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	519794	20.00	ng	0.00
39) Acenaphthene-d10	9.86	164	329732	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	555109	20.00	ng	0.00
76) Chrysene-d12	13.99	240	408879	20.00	ng	-0.01
86) Perylene-d12	15.45	264	423353	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	873193	94.65	ng	0.00
7) Phenol-d6	6.46	99	1083066	95.25	ng	0.00
23) Nitrobenzene-d5	7.39	82	587669	73.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	310301	101.57	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1291710	66.47	ng	0.00
79) Terphenyl-d14	12.95	244	1162153	63.25	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	203762	9.885	ng	99
71) Phenanthrene	11.37	178	82272	3.364	ng	98
75) Fluoranthene	12.56	202	213025	8.094	ng	99
78) Pyrene	12.79	202	217512	7.649	ng	98
81) Benzo(a)anthracene	13.98	228	102336	4.670	ng	96
83) Chrysene	14.02	228	98956	4.307	ng	97
87) Indeno(1,2,3-cd)pyrene	16.89	276	70355	3.038	ng	95
88) Benzo(b)fluoranthene	15.03	252	152684m	6.493	ng	
90) Benzo(a)pyrene	15.39	252	96239	4.679	ng	99
92) Benzo(g,h,i)perylene	17.32	276	69996	3.759	ng	98

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

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