

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060820\
 Data File : BM026287.D
 Acq On : 08 Jun 2020 15:19
 Operator : JU/CG
 Sample : L2839-20RE
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 NM21-COMP-2020-0-6RE

Manual Integrations
 APPROVED

mohammad
 6/9/2020 12:21:41 PM

Quant Time: Jun 08 15:56:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 15:00:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	244953	20.00	ng	0.00
21) Naphthalene-d8	10.20	136	1034275	20.00	ng	0.00
39) Acenaphthene-d10	14.09	164	615276	20.00	ng	0.00
64) Phenanthrene-d10	16.84	188	1271552	20.00	ng	0.00
76) Chrysene-d12	21.04	240	958434	20.00	ng	0.00
86) Perylene-d12	23.15	264	903612	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.08	112	1084084	78.25	ng	0.00
7) Phenol-d6	6.65	99	1643783	81.93	ng	0.00
23) Nitrobenzene-d5	8.59	82	1240562	58.89	ng	0.00
42) 2,4,6-Tribromophenol	15.59	330	625811	83.97	ng	0.00
45) 2-Fluorobiphenyl	12.70	172	2472213	60.99	ng	0.00
79) Terphenyl-d14	19.50	244	3346954	67.79	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.56	163	331371	7.363	ng	99
71) Phenanthrene	16.88	178	528740	7.985	ng	98
75) Fluoranthene	18.91	202	893909	11.775	ng	99
78) Pyrene	19.27	202	750986	11.859	ng	99
81) Benzo(a)anthracene	21.03	228	328991	5.719	ng	98
83) Chrysene	21.08	228	315304	5.751	ng	98
87) Indeno(1,2,3-cd)pyrene	25.20	276	180443	3.452	ng	96
88) Benzo(b)fluoranthene	22.53	252	379871	6.988	ng	98
89) Benzo(k)fluoranthene	22.57	252	141045m	2.671	ng	
90) Benzo(a)pyrene	23.06	252	270454	5.602	ng	99
92) Benzo(g,h,i)perylene	25.82	276	176340	4.247	ng	98

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

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