

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012519\
 Data File : BM018656.D
 Acq On : 25 Jan 2019 17:20
 Operator : JU/SJ
 Sample : K1157-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-03-011519

Manual Integrations
 APPROVED

mohammad
 1/29/2019 12:29:41 AM

Quant Time: Jan 28 02:43:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 22:01:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	225971	20.00	ng	0.00
21) Naphthalene-d8	10.52	136	893934	20.00	ng	0.00
39) Acenaphthene-d10	14.38	164	493557	20.00	ng	0.00
64) Phenanthrene-d10	17.13	188	1114244	20.00	ng	0.00
76) Chrysene-d12	21.33	240	1197653	20.00	ng	0.00
87) Perylene-d12	23.60	264	1245764	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	1162676	95.01	ng	0.00
7) Phenol-d6	6.90	99	1504757	96.81	ng	0.00
23) Nitrobenzene-d5	8.90	82	914656	59.31	ng	0.00
42) 2,4,6-Tribromophenol	15.87	330	618934	98.49	ng	-0.01
45) 2-Fluorobiphenyl	13.00	172	2089961	55.25	ng	0.00
79) Terphenyl-d14	19.76	244	3007347	52.93	ng	0.00
Target Compounds						
10) Phenol	6.93	94	77259	4.891	ng	94
50) Dimethylphthalate	13.84	163	169633	4.185	ng	99
71) Phenanthrene	17.17	178	890870	15.032	ng	99
72) Anthracene	17.26	178	217070	3.809	ng	98
75) Fluoranthene	19.19	202	970370	14.201	ng	98
78) Pyrene	19.56	202	809639	11.405	ng	99
81) Benzo(a)anthracene	21.31	228	479898	6.802	ng	99
83) Chrysene	21.36	228	410916	6.031	ng	96
86) Indeno(1,2,3-cd)pyrene	25.89	276	165510	2.107	ng	95
88) Benzo(b)fluoranthene	22.91	252	416354	5.887	ng	97
89) Benzo(k)fluoranthene	22.95	252	171601m	2.408	ng	
90) Benzo(a)pyrene	23.49	252	308375	4.554	ng	98
92) Benzo(a,h,i)perylene	26.60	276	145529	2.180	ng	97

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

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