

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011019\
 Data File : BM018526.D
 Acq On : 11 Jan 2019 01:41
 Operator : JU/SJ
 Sample : K1099-17
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-19-20

Manual Integrations
 APPROVED

Sohil
 1/11/2019 4:41:13 PM

Quant Time: Jan 11 04:19:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 10 05:57:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	784117	20.00	ng	0.00
21) Naphthalene-d8	10.55	136	3199203	20.00	ng	0.00
39) Acenaphthene-d10	14.40	164	1763445	20.00	ng	0.00
64) Phenanthrene-d10	17.16	188	3686766	20.00	ng	0.00
76) Chrysene-d12	21.35	240	2798830	20.00	ng	0.00
87) Perylene-d12	23.64	264	2635139	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	5986493	140.97	ng	0.00
7) Phenol-d6	6.95	99	7831312	145.20	ng	0.01
23) Nitrobenzene-d5	8.93	82	4885999	88.54	ng	0.00
42) 2,4,6-Tribromophenol	15.90	330	3146714	140.14	ng	0.00
45) 2-Fluorobiphenyl	13.03	172	10861779	80.37	ng	0.00
79) Terphenyl-d14	19.79	244	11594023	87.32	ng	0.00
Target Compounds						
10) Phenol	6.96	94	129220	2.358	ng	Qvalue # 70
50) Dimethylphthalate	13.87	163	2908420	20.082	ng	100
71) Phenanthrene	17.20	178	583776	2.977	ng	99
75) Fluoranthene	19.22	202	1282804	5.674	ng	98
78) Pyrene	19.58	202	1299739	7.834	ng	99
81) Benzo(a)anthracene	21.33	228	736623	4.468	ng	96
83) Chrysene	21.39	228	712621	4.476	ng	97
86) Indeno(1,2,3-cd)pyrene	25.94	276	609758	3.321	ng	# 91
88) Benzo(b)fluoranthene	22.94	252	1075422	7.189	ng	97
89) Benzo(k)fluoranthene	22.99	252	359456m	2.384	ng	
90) Benzo(a)pyrene	23.53	252	808326	5.644	ng	99
92) Benzo(a,h,i)perylene	26.66	276	602104	4.264	ng	98

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

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