

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
 Data File : BM020302.D
 Acq On : 12 May 2019 17:19
 Operator : JU/SJ
 Sample : K2768-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P026-SS015-0206-01

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:07:33 PM

Quant Time: May 13 07:01:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	55900	20.00	ng	-0.03
21) Naphthalene-d8	10.56	136	188356	20.00	ng	-0.03
39) Acenaphthene-d10	14.41	164	116352	20.00	ng	-0.02
64) Phenanthrene-d10	17.16	188	287412	20.00	ng	-0.02
76) Chrysene-d12	21.34	240	333420	20.00	ng	-0.02
87) Perylene-d12	23.63	264	399769	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	417081	121.96	ng	-0.02
7) Phenol-d6	6.93	99	545982	116.86	ng	-0.02
23) Nitrobenzene-d5	8.93	82	364911	76.49	ng	-0.03
42) 2,4,6-Tribromophenol	15.90	330	218954	121.24	ng	-0.02
45) 2-Fluorobiphenyl	13.03	172	611638	64.68	ng	-0.03
79) Terphenyl-d14	19.79	244	1127290	58.97	ng	-0.02
Target Compounds						
50) Dimethylphthalate	13.87	163	87715	8.870	ng	98
71) Phenanthrene	17.20	178	48933	3.084	ng	98
75) Fluoranthene	19.22	202	116236	5.790	ng	99
78) Pyrene	19.59	202	232757	10.398	ng	98
81) Benzo(a)anthracene	21.33	228	84403m	3.610	ng	
83) Chrysene	21.39	228	120942	5.333	ng	98
86) Indeno(1,2,3-cd)pyrene	25.94	276	77237	2.863	ng	# 100
88) Benzo(b)fluoranthene	22.94	252	149118m	5.649	ng	
90) Benzo(a)pyrene	23.53	252	92177	3.844	ng	96
92) Benzo(g,h,i)perylene	26.66	276	93352	3.943	ng	99

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

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