

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
 Data File : BM020296.D
 Acq On : 12 May 2019 13:44
 Operator : JU/SJ
 Sample : K2766-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P026-SS006-1824-01

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:07:25 PM

Quant Time: May 13 05:37:49 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	60352	20.00	ng	-0.03
21) Naphthalene-d8	10.56	136	200996	20.00	ng	-0.03
39) Acenaphthene-d10	14.41	164	119974	20.00	ng	-0.02
64) Phenanthrene-d10	17.16	188	293107	20.00	ng	-0.02
76) Chrysene-d12	21.35	240	324590	20.00	ng	-0.02
87) Perylene-d12	23.63	264	403320	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	462689	125.32	ng	-0.02
7) Phenol-d6	6.93	99	610022	120.94	ng	-0.02
23) Nitrobenzene-d5	8.93	82	413575	81.24	ng	-0.03
42) 2,4,6-Tribromophenol	15.90	330	242033	129.97	ng	-0.02
45) 2-Fluorobiphenyl	13.03	172	767320	78.69	ng	-0.03
79) Terphenyl-d14	19.79	244	1319741	70.92	ng	-0.02
Target Compounds						
50) Dimethylphthalate	13.87	163	67056	6.576	ng	98
75) Fluoranthene	19.22	202	54689	2.671	ng	100
78) Pyrene	19.59	202	104881	4.813	ng	99
83) Chrysene	21.39	228	64617	2.927	ng	96
88) Benzo(b)fluoranthene	22.94	252	78954m	2.965	ng	
92) Benzo(g,h,i)perylene	26.66	276	59933	2.509	ng	98

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

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