

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
 Data File : BM020301.D
 Acq On : 12 May 2019 16:43
 Operator : JU/SJ
 Sample : K2768-21
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P026-SS016-1824-01

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:07:32 PM

Quant Time: May 13 06:59:21 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	59959	20.00	ng	-0.03
21) Naphthalene-d8	10.56	136	197853	20.00	ng	-0.03
39) Acenaphthene-d10	14.41	164	116522	20.00	ng	-0.02
64) Phenanthrene-d10	17.16	188	271357	20.00	ng	-0.02
76) Chrysene-d12	21.34	240	294351	20.00	ng	-0.02
87) Perylene-d12	23.63	264	366727	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	436483	118.99	ng	-0.02
7) Phenol-d6	6.93	99	571327	114.01	ng	-0.02
23) Nitrobenzene-d5	8.93	82	394866	78.79	ng	-0.03
42) 2,4,6-Tribromophenol	15.90	330	218880	121.02	ng	-0.02
45) 2-Fluorobiphenyl	13.03	172	739740	78.11	ng	-0.03
79) Terphenyl-d14	19.79	244	1131557	67.05	ng	-0.02
Target Compounds						
50) Dimethylphthalate	13.87	163	125423	12.665	ng	99
71) Phenanthrene	17.20	178	30085	2.008	ng	96
75) Fluoranthene	19.22	202	59479	3.138	ng	99
78) Pyrene	19.59	202	106639	5.396	ng	99
81) Benzo(a)anthracene	21.33	228	44892m	2.175	ng	
83) Chrysene	21.38	228	53434m	2.669	ng	
88) Benzo(b)fluoranthene	22.94	252	49112m	2.028	ng	
90) Benzo(a)pyrene	23.53	252	44661	2.030	ng	# 94

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

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