

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
 Data File : BM019971.D
 Acq On : 19 Apr 2019 01:48
 Operator : JU/SJ
 Sample : K2442-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-20-S

Quant Time: Apr 19 02:23:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM041519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 19 00:56:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	118690	20.00	ng	0.00
21) Naphthalene-d8	10.29	136	448369	20.00	ng	0.00
39) Acenaphthene-d10	14.18	164	248912	20.00	ng	0.00
64) Phenanthrene-d10	16.94	188	545512	20.00	ng	0.00
76) Chrysene-d12	21.16	240	549191	20.00	ng	0.00
87) Perylene-d12	23.34	264	682322	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	615459	84.04	ng	0.00
7) Phenol-d6	6.71	99	832388	75.52	ng	0.00
23) Nitrobenzene-d5	8.69	82	523614	52.23	ng	0.00
42) 2,4,6-Tribromophenol	15.69	330	310895	77.59	ng	0.00
45) 2-Fluorobiphenyl	12.78	172	1037300	51.93	ng	0.00
79) Terphenyl-d14	19.59	244	1567260	50.34	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.66	163	123804	5.639	ng	97
75) Fluoranthene	19.02	202	126749	3.402	ng	99
78) Pyrene	19.39	202	125283	3.445	ng	99
81) Benzo(a)anthracene	21.14	228	82697	2.377	ng	97
83) Chrysene	21.20	228	73450	2.201	ng	96
88) Benzo(b)fluoranthene	22.70	252	107837	2.380	ng	# 94

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

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