

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031519\
 Data File : BM019182.D
 Acq On : 15 Mar 2019 12:33
 Operator : JU/SJ
 Sample : K1817-14
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PST-027-B116-030619

Manual Integrations
 APPROVED

Sohil
 3/18/2019 3:04:23 PM

Quant Time: Mar 16 02:04:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 17:30:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	206582	20.00	ng	0.00
21) Naphthalene-d8	10.44	136	955805	20.00	ng	-0.01
39) Acenaphthene-d10	14.31	164	616499	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1388090	20.00	ng	-0.01
76) Chrysene-d12	21.26	240	1347691	20.00	ng	-0.01
87) Perylene-d12	23.50	264	1225006	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	1225169	96.05	ng	0.00
7) Phenol-d6	6.83	99	1910124	102.37	ng	-0.01
23) Nitrobenzene-d5	8.82	82	1134528	56.11	ng	-0.01
42) 2,4,6-Tribromophenol	15.81	330	808238	88.79	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	2535386	53.49	ng	-0.01
79) Terphenyl-d14	19.70	244	3876002	57.10	ng	0.00
Target Compounds						
10) Phenol	6.86	94	42110	2.094	ng	86
50) Dimethylphthalate	13.77	163	206691	3.919	ng	99
71) Phenanthrene	17.10	178	222574	2.736	ng	99
75) Fluoranthene	19.13	202	339255	3.634	ng	99
78) Pyrene	19.49	202	422046	4.950	ng	99
81) Benzo(a)anthracene	21.24	228	170112	2.012	ng	98
83) Chrysene	21.30	228	201010	2.457	ng	97
88) Benzo(b)fluoranthene	22.83	252	199297m	2.511	ng	

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

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