

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101319\
 Data File : BP000685.D
 Acq On : 12 Oct 2019 22:23
 Operator : JU
 Sample : K5266-14RE
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AMENDED-COMPOSITE-C RE

Manual Integrations
 APPROVED

mohammad
 10/16/2019 8:08:10 AM

Quant Time: Oct 14 06:24:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	255775	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	909922	20.00	ng	-0.02
39) Acenaphthene-d10	9.79	164	511252	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	935337	20.00	ng	-0.02
76) Chrysene-d12	13.97	240	798958	20.00	ng	-0.01
87) Perylene-d12	15.45	264	882801	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	561292	35.27	ng	-0.02
7) Phenol-d6	6.39	99	875828	45.68	ng	-0.02
23) Nitrobenzene-d5	7.31	82	1377594	92.47	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	178468	32.76	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	2586484	81.86	ng	-0.02
79) Terphenyl-d14	12.90	244	3039522	77.03	ng	-0.01
Target Compounds						
10) Phenol	6.40	94	76097	3.876	ng	92
50) Dimethylphthalate	9.51	163	470822	13.254	ng	100
71) Phenanthrene	11.32	178	110572	2.354	ng	99
75) Fluoranthene	12.52	202	154385	2.926	ng	99
78) Pyrene	12.75	202	147704	2.680	ng	99
88) Benzo(b)fluoranthene	15.02	252	104068m	2.078	ng	

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

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