

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101019\
 Data File : BP000627.D
 Acq On : 10 Oct 2019 19:40
 Operator : JU
 Sample : K5285-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-2-S

Manual Integrations
 APPROVED

mohammad
 10/14/2019 8:40:28 AM

Quant Time: Oct 11 05:07:11 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	296881	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	1090581	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	590924	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	1047985	20.00	ng	-0.02
76) Chrysene-d12	13.96	240	856153	20.00	ng	-0.02
87) Perylene-d12	15.43	264	947396	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1422036	77.00	ng	-0.02
7) Phenol-d6	6.39	99	1826914	82.09	ng	-0.02
23) Nitrobenzene-d5	7.31	82	988802	55.38	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	469733	74.60	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1841301	50.42	ng	-0.02
79) Terphenyl-d14	12.90	244	1946550	46.03	ng	-0.02
Target Compounds						
10) Phenol	6.40	94	78199	3.432	ng	92
50) Dimethylphthalate	9.51	163	356135	8.673	ng	99
75) Fluoranthene	12.52	202	288834	4.886	ng	98
78) Pyrene	12.75	202	265009	4.488	ng	99
81) Benzo(a)anthracene	13.95	228	168807	3.232	ng	95
83) Chrysene	13.99	228	145658	2.841	ng	97
88) Benzo(b)fluoranthene	15.01	252	240307m	4.471	ng	
90) Benzo(a)pyrene	15.37	252	160249	3.198	ng	96
92) Benzo(g,h,i)perylene	17.33	276	124039	2.512	ng	99

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

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