

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092619\
 Data File : BP000420.D
 Acq On : 26 Sep 2019 14:20
 Operator : JU
 Sample : K5008-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TWP-5

Manual Integrations
 APPROVED

mohammad
 9/27/2019 1:46:50 PM

Quant Time: Sep 27 04:25:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	263034	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	957602	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	520991	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	906419	20.00	ng	0.00
76) Chrysene-d12	14.00	240	755507	20.00	ng	0.00
87) Perylene-d12	15.49	264	607855	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1394449	91.52	ng	0.00
7) Phenol-d6	6.42	99	1351442	72.36	ng	0.00
23) Nitrobenzene-d5	7.35	82	1231817	83.00	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	646868	125.18	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2514251	86.86	ng	0.00
79) Terphenyl-d14	12.94	244	2881142	80.00	ng	0.00
Target Compounds						
10) Phenol	6.43	94	69420	3.272	ng	92
37) 2-Methylnaphthalene	8.79	142	61049	2.014	ng	99
38) 1-Methylnaphthalene	8.88	142	132474m	4.665	ng	
50) Dimethylphthalate	9.55	163	225569	6.355	ng	# 98
52) Acenaphthene	9.86	154	128558	4.483	ng	98

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

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