

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091719\
 Data File : BP000275.D
 Acq On : 17 Sep 2019 16:13
 Operator : HP/JU
 Sample : K4918-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-8

Manual Integrations
 APPROVED

mohammad
 9/19/2019 4:46:14 PM

Quant Time: Oct 05 02:11:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 16 18:59:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	275520	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	969921	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	465678	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	306803	20.00	ng	0.04
76) Chrysene-d12	14.06	240	359561	20.00	ng	0.06
87) Perylene-d12	15.53	264	543541	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1552171	97.26	ng	0.01
7) Phenol-d6	6.43	99	1921535	98.22	ng	0.00
23) Nitrobenzene-d5	7.35	82	912936	60.73	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	396260	85.79	ng	0.02
45) 2-Fluorobiphenyl	9.15	172	2174517	84.04	ng	0.00
79) Terphenyl-d14	13.00	244	781775	45.61	ng	0.07
Target Compounds						
50) Dimethylphthalate	9.55	163	361407	11.391	ng	Qvalue # 95
71) Phenanthrene	11.39	178	85088	5.534	ng	# 7
88) Benzo(b)fluoranthene	15.10	252	84220m	2.597	ng	

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

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