

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000867.D
 Acq On : 18 Oct 2019 04:41
 Operator : JU
 Sample : K5392-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-22A-D

Quant Time: Oct 18 08:39:46 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	184567	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	643427	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	340874	20.00	ng	-0.02
64) Phenanthrene-d10	11.30	188	611476	20.00	ng	-0.01
76) Chrysene-d12	13.97	240	531089	20.00	ng	-0.01
87) Perylene-d12	15.45	264	575921	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	1027295	89.47	ng	-0.02
7) Phenol-d6	6.39	99	1287468	93.05	ng	-0.02
23) Nitrobenzene-d5	7.31	82	692863	65.77	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	371132	102.17	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1573960	74.71	ng	-0.02
79) Terphenyl-d14	12.90	244	1815116	69.20	ng	-0.01
Target Compounds						
10) Phenol	6.40	94	34546	2.439	ng	95
50) Dimethylphthalate	9.51	163	404091	17.061	ng	99

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

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