

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000893.D
 Acq On : 18 Oct 2019 22:16
 Operator : JU
 Sample : K5420-11
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-6

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:39:42 AM

Quant Time: Oct 19 05:54:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 16:40:44 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	157487	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	562847	20.00	ng	0.00
39) Acenaphthene-d10	9.79	164	310646	20.00	ng	0.00
64) Phenanthrene-d10	11.29	188	585258	20.00	ng	0.00
76) Chrysene-d12	13.96	240	528311	20.00	ng	0.00
87) Perylene-d12	15.44	264	606740	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	1162983	118.70	ng	0.00
7) Phenol-d6	6.39	99	1457144	123.42	ng	0.00
23) Nitrobenzene-d5	7.30	82	821872	89.18	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	493265	149.01	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1830639	95.35	ng	0.00
79) Terphenyl-d14	12.90	244	2213134	84.82	ng	0.00
Target Compounds						
10) Phenol	6.40	94	35028	2.898	ng	86
50) Dimethylphthalate	9.50	163	404894	18.758	ng	99
74) Di-n-butylphthalate	11.86	149	82399	2.504	ng	99

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

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