

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000885.D
 Acq On : 18 Oct 2019 19:02
 Operator : JU
 Sample : K5292-07RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20180657-AMENDED-COMP-DRE

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:39:29 AM

Quant Time: Oct 19 05:28:11 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	152560	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	538217	20.00	ng	0.00
39) Acenaphthene-d10	9.79	164	301229	20.00	ng	0.00
64) Phenanthrene-d10	11.29	188	549515	20.00	ng	0.00
76) Chrysene-d12	13.96	240	499353	20.00	ng	0.00
87) Perylene-d12	15.44	264	564611	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	442m	0.05	ng	0.04
7) Phenol-d6	6.39	99	6401	0.56	ng	0.00
23) Nitrobenzene-d5	7.30	82	753704	85.53	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	1757	0.55	ng	0.00
45) 2-Fluorobiphenyl	9.11	172	1776936	95.44	ng	0.00
79) Terphenyl-d14	12.90	244	2255366	91.45	ng	0.00
Target Compounds						
31) Naphthalene	8.05	128	63714	2.535	ng	99
50) Dimethylphthalate	9.50	163	48305	2.308	ng	99

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

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