

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101419\
 Data File : BP000716.D
 Acq On : 14 Oct 2019 17:31
 Operator : JU
 Sample : K5348-13RE
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20190881-COMPOSITE-MRE

Manual Integrations
 APPROVED

mohammad
 10/16/2019 8:11:46 AM

Quant Time: Oct 15 04:51:48 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	205535	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	731793	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	402323	20.00	ng	-0.02
64) Phenanthrene-d10	11.30	188	737034	20.00	ng	-0.01
76) Chrysene-d12	13.97	240	655743	20.00	ng	0.00
87) Perylene-d12	15.46	264	769218	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	73119	5.72	ng	-0.01
7) Phenol-d6	6.39	99	658378	42.73	ng	-0.02
23) Nitrobenzene-d5	7.31	82	792811	66.17	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	14721	3.43	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1496775	60.19	ng	-0.02
79) Terphenyl-d14	12.90	244	1742907	53.82	ng	-0.01
Target Compounds						
10) Phenol	6.40	94	63841	4.047	ng	92
50) Dimethylphthalate	9.51	163	269896	9.655	ng	99
75) Fluoranthene	12.53	202	148173	3.564	ng	99
78) Pyrene	12.76	202	154044	3.406	ng	99
81) Benzo(a)anthracene	13.96	228	94931	2.373	ng	# 82
83) Chrysene	14.00	228	87050	2.217	ng	97
84) Bis(2-ethylhexyl)phthalate	13.96	149	62962	2.540	ng	99
88) Benzo(b)fluoranthene	15.02	252	143459m	3.287	ng	
90) Benzo(a)pyrene	15.39	252	104327	2.565	ng	# 92

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

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