

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061419\
 Data File : BE100022.D
 Acq On : 14 Jun 2019 14:31
 Operator : JU/SJ
 Sample : K3284-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-TT-AOC22-MW03-20190609

Manual Integrations
 APPROVED

mohammad
 6/17/2019 4:51:33 PM

Quant Time: Jun 17 07:03:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 14:57:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	1007	0.40	ng	-0.01
7) Naphthalene-d8	10.59	136	3741	0.40	ng	-0.01
13) Acenaphthene-d10	14.46	164	3468m	0.40	ng	-0.01
19) Phenanthrene-d10	17.21	188	10132	0.40	ng	0.00
27) Chrysene-d12	21.40	240	11291	0.40	ng	0.00
34) Perylene-d12	23.94	264	12565	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.36	112	559	0.21	ng	0.01
5) Phenol-d6	6.95	99	370	0.11	ng	-0.01
8) Nitrobenzene-d5	8.95	82	1871	0.53	ng	-0.01
11) 2-Methylnaphthalene-d10	12.18	152	2879	0.44	ng	-0.03
14) 2,4,6-Tribromophenol	15.96	330	916	0.36	ng	-0.01
15) 2-Fluorobiphenyl	13.07	172	4653	0.35	ng	-0.03
25) Fluoranthene-d10	19.25	212	56950	0.39	ng	0.00
29) Terphenyl-d14	19.85	244	10665	0.52	ng	0.00
Target Compounds						
9) Naphthalene	10.64	128	10062	0.249	ng	# 82
16) Acenaphthylene	14.18	152	463	0.027	ng	# 1
17) Acenaphthene	14.51	154	757	0.081	ng	# 87

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

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