

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031319\
 Data File : BE098928.D
 Acq On : 14 Mar 2019 2:31
 Operator : JU/SJ
 Sample : K1856-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-TT-AOC22-MW07-20190311

Manual Integrations
 APPROVED

Sohil
 3/14/2019 11:09:29 AM

Quant Time: Mar 14 09:52:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	802	0.40	ng	0.01
7) Naphthalene-d8	10.60	136	3680	0.40	ng	0.01
13) Acenaphthene-d10	14.46	164	2601	0.40	ng	0.00
19) Phenanthrene-d10	17.21	188	6390	0.40	ng	0.00
27) Chrysene-d12	21.39	240	7682	0.40	ng	-0.01
34) Perylene-d12	23.89	264	7687	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.41	112	321	0.13	ng	0.02
5) Phenol-d6	7.08	99	179m	0.05	ng	0.09
8) Nitrobenzene-d5	9.02	82	1007	0.26	ng	0.05
11) 2-Methylnaphthalene-d10	12.21	152	1988	0.29	ng	0.01
14) 2,4,6-Tribromophenol	15.97	330	433	0.33	ng	0.00
15) 2-Fluorobiphenyl	13.09	172	4050	0.38	ng	0.00
25) Fluoranthene-d10	19.24	212	30599	0.30	ng	0.00
29) Terphenyl-d14	19.84	244	7834	0.42	ng	0.00
Target Compounds						
23) Phenanthrene	17.25	178	669	0.037	ng	Qvalue # 95
26) Fluoranthene	19.27	202	1740	0.068	ng	# 99
28) Pyrene	19.64	202	1457	0.062	ng	# 97
30) Benzo(a)anthracene	21.38	228	608	0.025	ng	# 72
31) Chrysene	21.44	228	958	0.037	ng	# 83
32) Bis(2-ethylhexyl)phthalate	21.29	149	2522	0.049	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.54	276	580	0.021	ng	# 73
35) Benzo(b)fluoranthene	23.14	252	1238	0.049	ng	# 46
36) Benzo(k)fluoranthene	23.18	252	545m	0.021	ng	
37) Benzo(a)pyrene	23.79	252	658	0.027	ng	# 13
39) Benzo(a,h,i)perylene	27.34	276	589	0.025	ng	# 60

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

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