

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE011018\
 Data File : BE095069.D
 Acq On : 10 Jan 2018 12:47
 Operator : SJ/JU
 Sample : J1041-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 OUTFALL001-180103

Quant Time: Jan 10 14:02:38 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 10 13:00:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	6041	0.40	ng	0.00
7) Naphthalene-d8	11.31	136	13105	0.40	ng	0.00
13) Acenaphthene-d10	15.05	164	6706	0.40	ng	0.00
19) Phenanthrene-d10	17.76	188	16624	0.40	ng	0.02
27) Chrysene-d12	21.85	240	18287	0.40	ng	0.01
34) Perylene-d12	24.51	264	14243	0.40	ng	0.02

System Monitoring Compounds						
4) 2-Fluorophenol	5.94	112	1830	0.20	ng	0.00
5) Phenol-d6	7.59	99	1698	0.12	ng	0.01
8) Nitrobenzene-d5	9.65	82	4607	0.43	ng	0.02
11) 2-Methylnaphthalene-d10	12.85	152	8767	0.41	ng	0.00
14) 2,4,6-Tribromophenol	16.52	330	609	0.39	ng	0.02
15) 2-Fluorobiphenyl	13.68	172	17380	0.58	ng	0.00
25) Fluoranthene-d10	19.73	212	93362	0.43	ng	0.00
29) Terphenyl-d14	20.30	244	23749	0.69	ng	0.01

Target Compounds					Qvalue	
9) Naphthalene	11.36	128	17587	0.114	ng	99
12) 2-Methylnaphthalene	12.92	142	3986	0.104	ng	96
16) Acenaphthylene	14.76	152	3709	0.103	ng	99
17) Acenaphthene	15.11	154	2739	0.115	ng	98
18) Fluorene	16.07	166	3453	0.117	ng	# 78
23) Phenanthrene	17.79	178	6655	0.128	ng	98
24) Anthracene	17.88	178	6491	0.122	ng	# 77
26) Fluoranthene	19.76	202	8025	0.126	ng	98
28) Pyrene	20.11	202	8133	0.131	ng	99
30) Benzo(a)anthracene	21.82	228	6448	0.121	ng	# 64
31) Chrysene	21.89	228	8636	0.144	ng	# 64
33) Indeno(1,2,3-cd)pyrene	27.36	276	5288	0.108	ng	95
35) Benzo(b)fluoranthene	23.68	252	6197	0.118	ng	# 94
36) Benzo(k)fluoranthene	23.74	252	5937	0.112	ng	# 88
37) Benzo(a)pyrene	24.39	252	4920	0.106	ng	# 84
38) Dibenzo(a,h)anthracene	27.38	278	4000	0.093	ng	94
39) Benzo(g,h,i)perylene	28.24	276	4980	0.116	ng	# 94

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

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