

Data Path : U:\HPCHEM1\BNA E\DATA\BE012518\
 Data File : BE095112.D
 Acq On : 25 Jan 2018 12:28
 Operator : SJ/JU
 Sample : J1236-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TW-2

Manual Integrations
 APPROVED

Sohil
 1/26/2018 11:21:55 AM

Quant Time: Jan 25 20:16:11 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE012318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 15:47:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	7962	0.40	nq	0.01
7) Naphthalene-d8	11.31	136	13198m	0.40	nq	0.00
13) Acenaphthene-d10	15.03	164	14489	0.40	nq	0.00
19) Phenanthrene-d10	17.74	188	15795	0.40	nq	0.00
27) Chrysene-d12	21.84	240	16947	0.40	nq	0.00
34) Perylene-d12	24.48	264	15993	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.97	112	3555	0.28	nq	0.02
5) Phenol-d6	7.58	99	12168m	0.62	nq	0.00
8) Nitrobenzene-d5	9.65	82	27702	2.31	nq	0.02
11) 2-Methylnaphthalene-d10	12.83	152	8270	0.38	nq	0.00
14) 2,4,6-Tribromophenol	16.50	330	7145	1.55	nq	0.00
15) 2-Fluorobiphenyl	13.68	172	128036	2.01	nq	0.00
25) Fluoranthene-d10	19.71	212	86683	0.44	nq	0.00
29) Terphenyl-d14	20.28	244	750541	21.64	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	11.36	128	15352444	101.655	nq	99
12) 2-Methylnaphthalene	12.91	142	207024	7.996	nq	97
16) Acenaphthylene	14.77	152	4661	0.060	nq	# 69
17) Acenaphthene	15.09	154	19017	0.377	nq	96
18) Fluorene	16.05	166	36831	0.584	nq	# 81
22) Pentachlorophenol	17.39	266	195m	0.136	nq	
23) Phenanthrene	17.77	178	201219	4.122	nq	98
24) Anthracene	17.86	178	42315	0.972	nq	96
26) Fluoranthene	19.74	202	242623	4.140	nq	97
28) Pvrene	20.10	202	208041	2.904	nq	# 85
30) Benzo(a)anthracene	21.81	228	101983	1.924	nq	98
31) Chrysene	21.87	228	87867	1.639	nq	97
32) Bis(2-ethylhexyl)phthalate	21.70	149	1476855	20.030	nq	100
33) Indeno(1,2,3-cd)pyrene	27.30	276	51458	1.056	nq	# 91
35) Benzo(b)fluoranthene	23.66	252	104277	1.838	nq	# 91
36) Benzo(k)fluoranthene	23.71	252	40950m	0.705	nq	
37) Benzo(a)pvrene	24.36	252	64314	1.237	nq	95
38) Dibenzo(a,h)anthracene	27.32	278	13232	0.302	nq	96
39) Benzo(g,h,i)perylene	28.18	276	51492	1.085	nq	# 92

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

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