

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040821\
 Data File : BE101264.D
 Acq On : 9 Apr 2021 6:21
 Operator : CG/JU
 Sample : M1966-03
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-110RR

Manual Integrations
 APPROVED

mohammad
 4/9/2021 11:19:37 AM

Quant Time: Apr 09 07:04:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-SIM-BE040621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 07 10:22:45 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.839	152	3023	0.40	ng	0.00
7) Naphthalene-d8	10.614	136	16101	0.40	ng	# 0.00
13) Acenaphthene-d10	14.442	164	15992	0.40	ng	0.00
19) Phenanthrene-d10	17.211	188	36882	0.40	ng	# 0.00
29) Chrysene-d12	21.365	240	30526	0.40	ng	# 0.01
36) Perylene-d12	23.841	264	25364	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.409	112	2523	0.35	ng	-0.01
5) Phenol-d6	6.998	99	1081	0.10	ng	0.00
8) Nitrobenzene-d5	8.977	82	5063	0.42	ng	0.00
11) 2-Methylnaphthalene-d10	12.195	152	9321m	0.25	ng	0.00
14) 2,4,6-Tribromophenol	15.956	330	1216	0.28	ng	0.00
15) 2-Fluorobiphenyl	13.074	172	14122	0.25	ng	0.00
27) Fluoranthene-d10	19.215	212	173393	0.27	ng	0.00
31) Terphenyl-d14	19.818	244	38830	0.47	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.370	88	936	0.20	ng	# 1
9) Naphthalene	10.666	128	1286776	7.35	ng	100
12) 2-Methylnaphthalene	12.267	142	4696	0.15	ng	# 91
16) Acenaphthylene	14.169	152	11241	0.18	ng	# 73
17) Acenaphthene	14.515	154	16868	0.36	ng	94
18) Fluorene	15.503	166	24795	0.40	ng	98
24) Pentachlorophenol	16.863	266	268	0.04	ng	# 77
25) Phenanthrene	17.241	178	81680m	0.75	ng	
26) Anthracene	17.332	178	61838	0.66	ng	98
28) Fluoranthene	19.249	202	78157	0.55	ng	98
30) Pyrene	19.605	202	64551	0.59	ng	98
32) Benzo(a)anthracene	21.341	228	22707	0.26	ng	# 91
33) Chrysene	21.401	228	19658	0.19	ng	# 90
34) Bis(2-ethylhexyl)phtha...	21.280	149	20342	0.33	ng	97
35) Indeno(1,2,3-cd)pyrene	26.448	276	11332	0.15	ng	99
37) Benzo(b)fluoranthene	23.078	252	17251	0.21	ng	# 87
38) Benzo(k)fluoranthene	23.128	252	13937	0.15	ng	# 84
39) Benzo(a)pyrene	23.727	252	12846	0.18	ng	# 84
40) Dibenzo(a,h)anthracene	26.477	278	7913	0.12	ng	# 85
41) Benzo(g,h,i)perylene	27.251	276	10281	0.14	ng	# 90

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

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