

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122416\
 Data File : BM008648.D
 Acq On : 25 Dec 2016 00:44
 Operator : UM/SJ
 Sample : H6069-09DL 2X
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8X9DL

Manual Integrations
 APPROVED

sohil
 12/26/2016 6:28:08 PM

Quant Time: Dec 26 02:19:27 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 26 00:55:42 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	430	0.40	ng/ul	-0.06
2) Naphthalene-d8	10.44	136	1302	0.40	ng/ul	-0.04
6) Acenaphthene-d10	14.30	164	792	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.06	188	1370	0.40	ng/ul	-0.02
16) Chrysene-d12	21.26	240	1470	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	1581m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	207	0.10	ng/ul	-0.02
14) Fluoranthene-d10	19.10	212	613m	0.16	ng/ul	0.00
Target Compounds						Ovalue
3) Naphthalene	10.49	128	2397	0.65	ng/ul	96
5) 2-Methylnaphthalene	12.13	142	748	0.30	ng/ul	99
7) Acenaphthylene	14.01	152	287	0.08	ng/ul#	60
8) Acenaphthene	14.35	153	96	0.04	ng/ul#	83
9) Fluorene	15.36	166	82	0.03	ng/ul#	79
12) Phenanthrene	17.09	178	1435	0.34	ng/ul#	95
13) Anthracene	17.19	178	501m	0.12	ng/ul	
15) Fluoranthene	19.12	202	2699	0.53	ng/ul	97
17) Pyrene	19.48	202	2604	0.46	ng/ul	95
18) Benzo(a)anthracene	21.24	228	1975	0.41	ng/ul#	89
19) Chrysene	21.29	228	1968	0.37	ng/ul#	90
21) Benzo(b)fluoranthene	22.82	252	3992m	0.64	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	1440m	0.24	ng/ul	
23) Benzo(a)pyrene	23.39	252	2623	0.45	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.74	276	2513	0.37	ng/ul#	91
26) Benzo(a,h,i)perylene	26.43	276	2351	0.40	ng/ul	94

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

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