

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080919\
 Data File : BM022018.D
 Acq On : 11 Aug 2019 05:22
 Operator : HP/JU
 Sample : K4028-05DL 10X
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BENQ7DL

Manual Integrations
 APPROVED

Sohil
 8/13/2019 8:34:27 AM

Quant Time: Aug 12 01:43:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 09 01:29:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	1135	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	4827	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.27	164	2702	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.03	188	6013	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	5965	0.40	ng/ul	-0.02
20) Perylene-d12	23.43	264	7621	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	177	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	506	0.02	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.45	128	333	0.024	ng/ul#	55
5) 2-Methylnaphthalene	12.09	142	207	0.023	ng/ul	91
7) Acenaphthylene	13.99	152	1057	0.084	ng/ul#	82
8) Acenaphthene	14.33	153	533	0.052	ng/ul	93
9) Fluorene	15.33	166	649	0.057	ng/ul#	94
12) Phenanthrene	17.07	178	14587	0.834	ng/ul	96
13) Anthracene	17.16	178	3822	0.256	ng/ul	99
15) Fluoranthene	19.09	202	29023	1.238	ng/ul	97
17) Pyrene	19.45	202	35718	1.457	ng/ul	99
18) Benzo(a)anthracene	21.21	228	21253	0.996	ng/ul	98
19) Chrysene	21.26	228	21711	0.791	ng/ul	97
21) Benzo(b)fluoranthene	22.77	252	21271m	0.828	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	7934m	0.273	ng/ul	
23) Benzo(a)pyrene	23.33	252	17878	0.684	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.64	276	10073	0.323	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.64	278	3013	0.119	ng/ul#	68
26) Benzo(g,h,i)perylene	26.32	276	10547	0.375	ng/ul	97

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

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