

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021927.D
 Acq On : 08 Aug 2019 11:18
 Operator : HP/JU
 Sample : K4029-15DL 5X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BENY5DL

Manual Integrations
 APPROVED

mohammad
 8/9/2019 6:12:16 AM

Quant Time: Aug 08 13:58:01 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 : Thu Aug 08 06:54:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	707	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	3049	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.28	164	1677	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	3661	0.40	ng/ul	-0.02
16) Chrysene-d12	21.24	240	3428	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	5058	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	293	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	708	0.06	ng/ul	-0.01
Target Compounds				Ovalue		
7) Acenaphthylene	14.00	152	412	0.053	ng/ul#	70
12) Phenanthrene	17.07	178	2639	0.248	ng/ul	98
13) Anthracene	17.17	178	713	0.078	ng/ul#	74
15) Fluoranthene	19.10	202	6896	0.483	ng/ul	100
17) Pyrene	19.46	202	6759	0.480	ng/ul	95
18) Benzo(a)anthracene	21.22	228	3983	0.325	ng/ul	94
19) Chrysene	21.28	228	3870	0.245	ng/ul	95
21) Benzo(b)fluoranthene	22.79	252	5298m	0.311	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	2504m	0.130	ng/ul	
23) Benzo(a)pyrene	23.36	252	3537	0.204	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.68	276	2605	0.126	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.68	278	613	0.037	ng/ul#	23
26) Benzo(g,h,i)perylene	26.36	276	2197	0.118	ng/ul#	86

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

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