

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081219\
 Data File : BM022024.D
 Acq On : 12 Aug 2019 10:24
 Operator : HP/JU
 Sample : K4028-06DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BENY8DL

Manual Integrations
 APPROVED

Sohil
 8/13/2019 9:57:00 AM

Quant Time: Aug 12 14:19:01 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 : Mon Aug 12 14:06:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	745	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	3099	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.27	164	1684	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.02	188	3778	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	3915	0.40	ng/ul	-0.02
20) Perylene-d12	23.43	264	4227	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	232	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	680	0.05	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	641	0.073	ng/ul#	81
5) 2-Methylnaphthalene	12.09	142	337	0.058	ng/ul	91
7) Acenaphthylene	13.99	152	2460	0.313	ng/ul#	84
8) Acenaphthene	14.33	153	1334	0.208	ng/ul	98
9) Fluorene	15.33	166	1598	0.224	ng/ul	99
12) Phenanthrene	17.06	178	40122	3.650	ng/ul	97
13) Anthracene	17.16	178	7312	0.780	ng/ul	98
15) Fluoranthene	19.08	202	120026	8.151	ng/ul	96
17) Pyrene	19.45	202	110971	6.898	ng/ul	98
18) Benzo(a)anthracene	21.21	228	56060	4.003	ng/ul	99
19) Chrysene	21.26	228	66866	3.713	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	92275m	6.474	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	35382m	2.193	ng/ul	
23) Benzo(a)pyrene	23.33	252	61490	4.242	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.64	276	45806	2.646	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.64	278	10993	0.786	ng/ul	95
26) Benzo(a,h,i)perylene	26.32	276	46762	3.001	ng/ul	96

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

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