

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042718\
 Data File : BE096207.D
 Acq On : 27 Apr 2018 14:34
 Operator : SJ/JU
 Sample : J2449-03DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD8DL

Manual Integrations
 APPROVED

Sohil
 4/30/2018 6:58:46 PM

Quant Time: Apr 28 00:51:46 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 27 23:57:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	750	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2805	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1587	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3865	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	4822	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	4118	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	427	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	1242	0.07	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.23	128	1963	0.251	ng/ul#	90
5) 2-Methylnaphthalene	12.81	142	1082	0.207	ng/ul	99
7) Acenaphthylene	14.67	152	961	0.117	ng/ul#	87
8) Acenaphthene	15.01	153	768	0.130	ng/ul#	92
9) Fluorene	15.97	166	1464	0.188	ng/ul	94
12) Phenanthrene	17.70	178	39551	3.476	ng/ul#	89
13) Anthracene	17.78	178	6665	0.584	ng/ul#	90
15) Fluoranthene	19.67	202	69176	3.977	ng/ul	84
17) Pyrene	20.02	202	30506	3.154	ng/ul#	77
18) Benzo(a)anthracene	21.73	228	15465	1.059	ng/ul#	86
19) Chrysene	21.79	228	23960	1.786	ng/ul	98
21) Benzo(b)fluoranthene	23.55	252	27201m	1.744	ng/ul	
22) Benzo(k)fluoranthene	23.59	252	9414m	0.632	ng/ul	
23) Benzo(a)pyrene	24.23	252	19469	1.437	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.13	276	12373	0.678	ng/ul#	75
25) Dibenzo(a,h)anthracene	27.13	278	3027	0.198	ng/ul	98
26) Benzo(a,h,i)perylene	28.00	276	10803	0.680	ng/ul#	94

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

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