

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097297.D
 Acq On : 17 Aug 2018 15:08
 Operator : SJ/JU
 Sample : J4424-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YQ1

Quant Time: Aug 17 23:21:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 14:15:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	976	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	4907	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	2539	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	6336	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	6276	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	5604	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2520	0.32	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	7322	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	3293	0.275	ng/ul#	94
5) 2-Methylnaphthalene	11.80	142	1958	0.254	ng/ul	98
7) Acenaphthylene	13.73	152	3652	0.319	ng/ul	97
9) Fluorene	15.08	166	3280	0.365	ng/ul	93
12) Phenanthrene	16.81	178	10755	0.657	ng/ul#	90
13) Anthracene	16.90	178	1654	0.121	ng/ul#	93
15) Fluoranthene	18.84	202	8409	0.483	ng/ul	84
17) Pyrene	19.20	202	9411	0.496	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	10322	0.663	ng/ul#	85
19) Chrysene	21.01	228	7439	0.401	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	4334	0.286	ng/ul	87
22) Benzo(k)fluoranthene	22.58	252	9047	0.480	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.52	276	6759	0.388	ng/ul#	78
25) Dibenzo(a,h)anthracene	25.54	278	6147	0.436	ng/ul#	91
26) Benzo(g,h,i)perylene	26.22	276	8667	0.621	ng/ul#	92

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

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