

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096185.D
 Acq On : 26 Apr 2018 18:44
 Operator : SJ/JU
 Sample : J2431-14RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Apr 27 05:24:35 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 04:45:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	1033	0.40	ng/ul	0.00
2) Naphthalene-d8	11.19	136	3850	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	11680	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.67	188	4639	0.40	ng/ul	0.01
16) Chrysene-d12	21.78	240	5504m	0.40	ng/ul	0.04
20) Perylene-d12	24.38	264	5283	0.40	ng/ul	0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1309	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	3263m	0.15	ng/ul	0.05
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	2565011	239.272	ng/ul#	87
5) 2-Methylnaphthalene	12.84	142	2964507	412.746	ng/ul	99
7) Acenaphthylene	14.68	152	145530	2.417	ng/ul#	93
8) Acenaphthene	15.05	153	4835892	111.392	ng/ul	96
9) Fluorene	16.02	166	4500106	78.342	ng/ul#	90
11) Pentachlorophenol	17.36	266	845	0.757	ng/ul	89
12) Phenanthrene	17.75	178	14435003	1056.874	ng/ul#	88
13) Anthracene	17.85	178	3767087m	274.916	ng/ul	
15) Fluoranthene	19.72	202	12037483	576.589	ng/ul	84
17) Pyrene	20.08	202	10587400	958.840	ng/ul#	79
18) Benzo(a)anthracene	21.77	228	3955838	237.397	ng/ul#	84
19) Chrysene	21.82	228	2207089	144.141	ng/ul	93
21) Benzo(b)fluoranthene	23.59	252	1786142m	89.289	ng/ul	
22) Benzo(k)fluoranthene	23.64	252	547224m	28.615	ng/ul	
23) Benzo(a)pyrene	24.28	252	1136870	65.417	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	27.18	276	412194	17.603	ng/ul#	72
25) Dibenzo(a,h)anthracene	27.17	278	119649m	6.102	ng/ul	
26) Benzo(g,h,i)perylene	28.05	276	335001	16.435	ng/ul#	90

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

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