

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096084.D
 Acq On : 23 Apr 2018 14:35
 Operator : SJ/JU
 Sample : J2355-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Y26

Manual Integrations
 APPROVED

Sohil
 4/24/2018 7:09:53 PM

Quant Time: Apr 24 06:37:21 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 24 05:21:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1867	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5058	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2914	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6501m	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	6119m	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6850	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	1456	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	4208m	0.16	ng/ul	0.00
Target Compounds						
					Ovalue	
7) Acenaphthylene	14.70	152	313	0.021	ng/ul#	65
8) Acenaphthene	15.02	153	228	0.021	ng/ul#	85
9) Fluorene	15.99	166	408	0.033	ng/ul#	89
11) Pentachlorophenol	17.33	266	112m	0.081	ng/ul	
12) Phenanthrene	17.71	178	1396m	0.073	ng/ul	
13) Anthracene	17.80	178	1099m	0.058	ng/ul	
15) Fluoranthene	19.68	202	2646m	0.101	ng/ul	
17) Pyrene	20.04	202	3187	0.153	ng/ul#	84
18) Benzo(a)anthracene	21.75	228	4131m	0.167	ng/ul	
19) Chrysene	21.80	228	2422	0.136	ng/ul	91
21) Benzo(b)fluoranthene	23.57	252	2748	0.124	ng/ul	97
22) Benzo(k)fluoranthene	23.62	252	2385	0.108	ng/ul#	92
23) Benzo(a)pyrene	24.26	252	2632	0.123	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	27.17	276	2823	0.119	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.18	278	2148	0.112	ng/ul#	93
26) Benzo(a,h,i)perylene	28.04	276	2336	0.118	ng/ul#	91

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

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