

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062116\
 Data File : BM005892.D
 Acq On : 21 Jun 2016 14:04
 Operator : UM/SJ
 Sample : PB91207BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK07

Quant Time: Jun 21 14:35:39 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 21 08:19:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	167941	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	876052	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.32	164	594404	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.07	188	1504117	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	1810583	20.00	ng/ul	0.00
83) Perylene-d12	23.49	264	1842370	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.20	96	15052	3.51	ng/uL	0.00
5) Phenol-d5	6.85	99	608261	32.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.02	67	402214	37.50	ng/ul	0.00
9) 2-Chlorophenol-d4	7.21	132	457612	33.50	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	529458	34.18	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	239305	37.69	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	252366	39.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	499163	33.08	ng/ul	0.00
29) 4-Chloroaniline-d4	10.60	131	607416	43.00	ng/ul	0.00
43) Dimethylphthalate-d6	13.74	166	2140900	38.01	ng/ul	0.00
46) Acenaphthylene-d8	14.02	160	2377149	35.74	ng/ul	0.00
51) 4-Nitrophenol-d4	14.52	143	355211	36.56	ng/ul	0.00
57) Fluorene-d10	15.32	176	1739779	37.35	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.43	200	212684	35.09	ng/ul	0.00
70) Anthracene-d10	17.17	188	2863614	37.55	ng/ul	0.00
76) Pyrene-d10	19.47	212	3313048	34.26	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.36	264	3503097	37.32	ng/ul	0.00

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

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