

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011315\
 Data File : BM000166.D
 Acq On : 13 Jan 2015 13:30
 Operator : TP/IZ
 Sample : PB81299BL
 Misc : SBLK93
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK93

Quant Time: Jan 14 13:02:07 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 14 12:47:10 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	110035	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.62	136	447257	20.00	ng/ul	-0.02
35) Acenaphthene-d10	14.47	164	312877	20.00	ng/ul	-0.01
61) Phenanthrene-d10	17.21	188	634541	20.00	ng/ul	-0.02
75) Chrysene-d12	21.40	240	702263	20.00	ng/ul	-0.01
83) Perylene-d12	23.69	264	657353	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.32	96	14765	7.78	ng/uL	-0.01
5) Phenol-d5	6.99	99	341246	36.87	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	7.16	67	227322	38.85	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.36	132	256665	37.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	257979	35.03	ng/ul	-0.01
19) Nitrobenzene-d5	8.99	128	125599	40.61	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.70	143	127919	40.51	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.24	165	255654	36.65	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.76	131	349424	44.15	ng/ul	-0.02
43) Dimethylphthalate-d6	13.87	166	868191	37.62	ng/ul	-0.01
46) Acenaphthylene-d8	14.16	160	1056003	38.20	ng/ul	-0.02
51) 4-Nitrophenol-d4	14.65	143	103709	27.62	ng/ul	-0.01
57) Fluorene-d10	15.46	176	755449	38.12	ng/ul	-0.02
62) 4,6-Dinitro-2-methylphenol	15.57	200	97972	32.58	ng/ul	-0.02
70) Anthracene-d10	17.30	188	1172544	39.96	ng/ul	-0.01
76) Pyrene-d10	19.60	212	1292560	39.90	ng/ul	-0.01
87) Benzo(a)pyrene-d12	23.54	264	1247143	41.80	ng/ul	-0.02

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

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