

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\
 Data File : BM023731.D
 Acq On : 14 Nov 2019 23:53
 Operator : JU
 Sample : PB124582BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK82

Quant Time: Nov 15 07:51:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 15 06:30:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	396075	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1638358	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.19	164	1029770	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.94	188	2235656	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	2239506	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	2582287	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	64630	6.73	ng/uL	0.00
5) Phenol-d5	6.72	99	1210342	33.52	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	705855	34.02	ng/ul	0.00
9) 2-Chlorophenol-d4	7.07	132	962502	36.80	ng/ul	0.00
13) 4-Methylphenol-d8	8.25	113	962904	33.43	ng/ul	0.00
19) Nitrobenzene-d5	8.68	128	471508	40.21	ng/ul	0.00
22) 2-Nitrophenol-d4	9.40	143	545950	44.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	972522	34.60	ng/ul	0.00
29) 4-Chloroaniline-d4	10.45	131	1303787	42.24	ng/ul	0.00
44) Dimethylphthalate-d6	13.61	166	3022327	37.63	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	3458884	37.74	ng/ul	0.00
52) 4-Nitrophenol-d4	14.41	143	482999	35.23	ng/ul	0.00
58) Fluorene-d10	15.19	176	2562769	36.14	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	431345	33.35	ng/ul	0.00
71) Anthracene-d10	17.04	188	3957239	37.31	ng/ul	0.00
79) Pyrene-d10	19.34	212	4493047	38.48	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	5257672	38.69	ng/ul	0.00

Target Compounds	Ovalue

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

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