

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
 Data File : BM023918.D
 Acq On : 27 Nov 2019 18:33
 Operator : JU
 Sample : PB124947BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK47

Quant Time: Nov 28 00:21:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 27 17:16:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	470628	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	1981546	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.15	164	1218585	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	2405293	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	2089347	20.00	ng/ul	0.00
86) Perylene-d12	23.26	264	2462365	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.03	96	70021	6.62	ng/uL	0.00
5) Phenol-d5	6.68	99	1318279	31.96	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.84	67	806927	33.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.03	132	1062263	33.32	ng/ul	0.00
13) 4-Methylphenol-d8	8.21	113	1063828	32.69	ng/ul	0.00
19) Nitrobenzene-d5	8.65	128	529016	34.40	ng/ul	0.00
22) 2-Nitrophenol-d4	9.36	143	616582	35.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.90	165	1061398	32.06	ng/ul	0.00
29) 4-Chloroaniline-d4	10.41	131	1449913	39.91	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	3415380	36.07	ng/ul	0.00
47) Acenaphthylene-d8	13.84	160	3900421	35.15	ng/ul	0.00
52) 4-Nitrophenol-d4	14.38	143	491454	30.87	ng/ul	0.00
58) Fluorene-d10	15.15	176	2783812	34.94	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	459587	30.44	ng/ul	0.00
71) Anthracene-d10	17.00	188	3978492	34.63	ng/ul	0.00
79) Pyrene-d10	19.31	212	4200944	38.91	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.13	264	4732214	36.36	ng/ul	0.00

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

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