

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060120\
 Data File : BM026201.D
 Acq On : 01 Jun 2020 17:01
 Operator : JU/CG
 Sample : PB129218BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK18

Quant Time: Jun 01 17:39:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 01 16:59:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	108583	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	415135	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.12	164	241750	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.87	188	562158	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	617967	20.00	ng/ul	0.00
86) Perylene-d12	23.19	264	670010	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.05	96	24668	6.85	ng/uL	0.00
5) Phenol-d5	6.67	99	383523	38.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.82	67	257478	37.79	ng/ul	0.00
9) 2-Chlorophenol-d4	7.01	132	289663	37.57	ng/ul	0.00
13) 4-Methylphenol-d8	8.19	113	283911	38.25	ng/ul	0.00
19) Nitrobenzene-d5	8.62	128	128165	37.87	ng/ul	0.00
22) 2-Nitrophenol-d4	9.33	143	135131	40.36	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.87	165	247123	36.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.38	131	359272	52.70	ng/ul	0.00
44) Dimethylphthalate-d6	13.55	166	776269	40.04	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	882215	38.41	ng/ul	0.00
52) 4-Nitrophenol-d4	14.35	143	123487	35.97	ng/ul	0.00
58) Fluorene-d10	15.12	176	673781	39.96	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.25	200	100836	30.76	ng/ul	0.00
71) Anthracene-d10	16.97	188	1058839	38.22	ng/ul	0.00
79) Pyrene-d10	19.27	212	1218201	38.38	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.06	264	1382641	38.77	ng/ul	0.00

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

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