

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041421\
 Data File : BM029462.D
 Acq On : 14 Apr 2021 10:01
 Operator : CG/JU
 Sample : PB135558BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK558

Quant Time: Apr 14 10:34:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM041221.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 14 10:06:38 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	81112	20.00	ng/ul	0.00
20) Naphthalene-d8	10.46	136	327038	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.32	164	189011	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.06	188	342612	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	296578	20.00	ng/ul	0.00
87) Perylene-d12	23.44	264	318365	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	16343	7.31	ng/uL	0.00
4) Pyridine-d5	3.59	84	211157	35.62	ng/ul	0.00
7) Phenol-d5	6.85	99	275504	37.10	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.02	67	188575	38.91	ng/ul	0.00
11) 2-Chlorophenol-d4	7.21	132	219900	39.53	ng/ul	0.00
15) 4-Methylphenol-d8	8.38	113	213091	37.51	ng/ul	0.00
21) Nitrobenzene-d5	8.83	128	94214	37.21	ng/ul	0.00
24) 2-Nitrophenol-d4	9.55	143	106370	38.58	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.08	165	189377	36.53	ng/ul	0.00
31) 4-Chloroaniline-d4	10.59	131	264885	35.91	ng/ul	0.00
46) Dimethylphthalate-d6	13.73	166	571890	39.25	ng/ul	0.00
49) Acenaphthylene-d8	14.01	160	702196	40.39	ng/ul	0.00
54) 4-Nitrophenol-d4	14.52	143	86311	30.96	ng/ul	0.00
60) Fluorene-d10	15.31	176	461105	38.44	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.43	200	71342	32.10	ng/ul	0.00
73) Anthracene-d10	17.16	188	699447	42.25	ng/ul	0.00
80) Pyrene-d10	19.45	212	727157	46.26	ng/ul	0.00
91) Benzo(a)pyrene-d12	23.30	264	775824	44.68	ng/ul	0.00

Target Compounds Ovalue

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

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