

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009922.D
 Acq On : 18 Apr 2022 18:55
 Operator : CG/JU
 Sample : PB144146BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK146

Quant Time: Apr 19 00:57:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	121276	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	538891	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.587	164	387071	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.334	188	872975	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.416	240	862964	20.000	ng/u1	# 0.00
88) Perylene-d12	23.874	264	766540	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	17766	6.538	ng/uL	0.00
4) Pyridine-d5	3.793	84	235462	31.428	ng/u1	0.00
7) Phenol-d5	7.105	99	309442	29.063	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	207973	32.810	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	251954	31.371	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	268817	30.244	ng/u1	-0.01
21) Nitrobenzene-d5	9.110	128	134416	34.589	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	144071	33.815	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	260407	29.209	ng/u1	0.00
31) 4-Chloroaniline-d4	10.887	131	396500	31.225	ng/u1	0.00
46) Dimethylphthalate-d6	13.993	166	993066	32.661	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	1093682	30.362	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.763	143	147220	26.699	ng/u1	0.00
60) Fluorene-d10	15.575	176	846321	33.088	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	137474	26.583	ng/u1	0.00
73) Anthracene-d10	17.433	188	1386664	33.176	ng/u1	0.00
81) Pyrene-d10	19.663	212	1680760	35.661	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.716	264	1368609	34.285	ng/u1	-0.01

Target Compounds	Qvalue

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

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